LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100618\
Data File : VU027485.D
Acq On : 06 Oct 2018 03:41
Operator : MD/SY
Sample : J5268-04 10X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-11

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\82U100118W.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	1.759	172	182	207	rVB	357739	473647	37.12%	3.210%
2	1.948	232	241	259	rVB	199402	272790	21.38%	1.849%
3	2.051	266	273	285	rVB	14902	20058	1.57%	0.136%
4	2.186	307	315	327	rVB	18235	27675	2.17%	0.188%
5	2.305	341	352	368	rVB3	20859	42191	3.31%	0.286%
6	2.646	444	458	466	rBV6	6661	16759	1.31%	0.114%
7	2.710	466	478	480	rBV2	66819	97142	7.61%	0.658%
8	2.746	481	489	499	rVB	164940	303106	23.76%	2.054%
9	3.000	552	568	587	rBV	124391	263478	20.65%	1.785%
10	3.193	616	628	645	rVB4	18429	44195	3.46%	0.299%
11	3.305	648	663	686	rBV	84965	183057	14.35%	1.240%
12	3.427	689	701	713	rVV3	11793	25974	2.04%	0.176%
13	3.501	713	724	732	rVV2	15522	34124	2.67%	0.231%
14	3.559	733	742	757	rVB	30088	65670	5.15%	0.445%
15	3.659	759	773	787	rBV3	16689	40020	3.14%	0.271%
16	3.746	787	800	815	rVB9	8019	23739	1.86%	0.161%
17	3.894	827	846	862	rVB3	24907	64316	5.04%	0.436%
18	4.000	865	879	892	rVV2	12126	31213	2.45%	0.212%
19	4.096	892	909	928	rVV2	144240	387917	30.40%	2.629%
20	4.193	929	939	962	rVB7	9292	27553	2.16%	0.187%
21	4.736	1086	1108	1111	rBV10	6864	16013	1.26%	0.109%
22	4.784	1112	1123	1138	rVV2	24764	58195	4.56%	0.394%
23	4.887	1138	1155	1171	rVV	106779	246149	19.29%	1.668%
24	4.990	1171	1187	1204	rVV4	252055	633673	49.67%	4.294%
25	5.083	1208	1216	1241	rVB2	35625	96157	7.54%	0.652%
26	5.221	1241	1259	1275	rBV2	40218	102798	8.06%	0.697%
27	5.311	1275	1287	1312	rVB	132600	279837	21.93%	1.896%
28	5.485	1328	1341	1348	rBV2	26858	56627	4.44%	0.384%
29	5.546	1349	1360	1373	rVV4	42346	118086	9.26%	0.800%
30	5.623	1374	1384	1399	rVB3	21340	47938	3.76%	0.325%
31	5.784	1420	1434	1445	rBV2	23876	49248	3.86%	0.334%
32	5.890	1454	1467	1478	rBV	224931	440890	34.56%	2.988%
33	5.948	1479	1485	1511	rVB6	33170	97427	7.64%	0.660%
34	6.122	1529	1539	1552	rVB4	6706	12816	1.00%	0.087%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100618\
Data File : VU027485.D
Acq On : 06 Oct 2018 03:41
Operator : MD/SY
Sample : J5268-04 10X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-11

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\82U100118W.M
Title : SW846 8260

35	6.225	1559	1571	1589	rBV6	14451	32635	2.56%	0.221%
36	6.418	1611	1631	1644	rBV3	83991	205562	16.11%	1.393%
37	6.643	1691	1701	1715	rVB3	13174	25524	2.00%	0.173%
38	6.842	1744	1763	1773	rBV10	7163	22951	1.80%	0.156%
39	6.926	1776	1789	1808	rVB3	16786	37785	2.96%	0.256%
40	7.061	1809	1831	1842	rBV5	30521	83949	6.58%	0.569%
41	7.434	1936	1947	1964	rVB5	16705	34342	2.69%	0.233%
42	7.569	1964	1989	2008	rBV	438057	820556	64.31%	5.560%
43	8.048	2127	2138	2146	rBV2	7878	13431	1.05%	0.091%
44	9.089	2451	2462	2484	rBV	336854	573683	44.96%	3.887%
45	9.250	2500	2512	2536	rVB	416549	667970	52.35%	4.526%
46	9.376	2536	2551	2571	rBV	779903	1275892	100.00%	8.646%
47	9.781	2664	2677	2699	rVB	236575	377523	29.59%	2.558%
48	10.167	2786	2797	2808	rBV2	36384	55714	4.37%	0.378%
49	10.308	2828	2841	2869	rBV	333509	528072	41.39%	3.578%
50	10.588	2917	2928	2945	rBV	102369	160624	12.59%	1.088%
51	10.681	2945	2957	2962	rBV	264613	426429	33.42%	2.890%
52	10.771	2976	2985	3006	rVB	156165	239073	18.74%	1.620%
53	10.970	3035	3047	3063	rBV	185460	291220	22.82%	1.973%
54	11.147	3088	3102	3119	rBV	774080	1180984	92.56%	8.003%
55	11.430	3179	3190	3195	rBV	11017	15891	1.25%	0.108%
56	11.485	3195	3207	3223	rVV	436062	683311	53.56%	4.630%
57	11.572	3223	3234	3247	rVB	178252	271844	21.31%	1.842%
58	11.768	3283	3295	3304	rBV	197473	345576	27.09%	2.342%
59	11.810	3304	3308	3317	rVV	49567	66141	5.18%	0.448%
60	11.874	3317	3328	3347	rVB5	73500	146181	11.46%	0.991%
61	11.980	3353	3361	3371	rBV4	8875	13207	1.04%	0.089%
62	12.057	3374	3385	3399	rVB	18286	27905	2.19%	0.189%
63	12.147	3403	3413	3418	rBV	47252	74106	5.81%	0.502%
64	12.176	3418	3422	3433	rVV	38102	52285	4.10%	0.354%
65	12.250	3433	3445	3452	rVV	75783	116083	9.10%	0.787%
66	12.285	3453	3456	3466	rVB2	23505	29018	2.27%	0.197%
67	12.353	3466	3477	3500	rBV	65533	115850	9.08%	0.785%
68	12.552	3528	3539	3549	rVB2	16188	25038	1.96%	0.170%
69	12.652	3554	3570	3577	rBV	42299	65102	5.10%	0.441%
70	12.703	3577	3586	3597	rVB	62172	91730	7.19%	0.622%
71	12.964	3658	3667	3681	rVB	65631	101233	7.93%	0.686%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100618\
Data File : VU027485.D
Acq On : 06 Oct 2018 03:41
Operator : MD/SY
Sample : J5268-04 10X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-11

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\82U100118W.M
Title : SW846 8260

72	13.125	3696	3717	3729	rBV2	123063	213870	16.76%	1.449%
73	13.183	3731	3735	3745	rVV6	9324	14980	1.17%	0.102%
74	13.292	3760	3769	3786	rVB4	15923	27033	2.12%	0.183%
75	13.456	3812	3820	3829	rBV2	14617	21772	1.71%	0.148%
76	13.511	3829	3837	3844	rBV5	16670	27789	2.18%	0.188%
77	13.742	3898	3909	3929	rBV	191243	305482	23.94%	2.070%
78	13.861	3937	3946	3959	rVB2	8284	13397	1.05%	0.091%
79	14.154	4027	4037	4048	rBV3	9734	14333	1.12%	0.097%
80	14.334	4084	4093	4106	rBV4	6459	14188	1.11%	0.096%
81	14.880	4253	4263	4279	rBV	35062	63715	4.99%	0.432%
82	15.063	4310	4320	4339	rBV2	26389	45815	3.59%	0.310%

Sum of corrected areas: 14757272

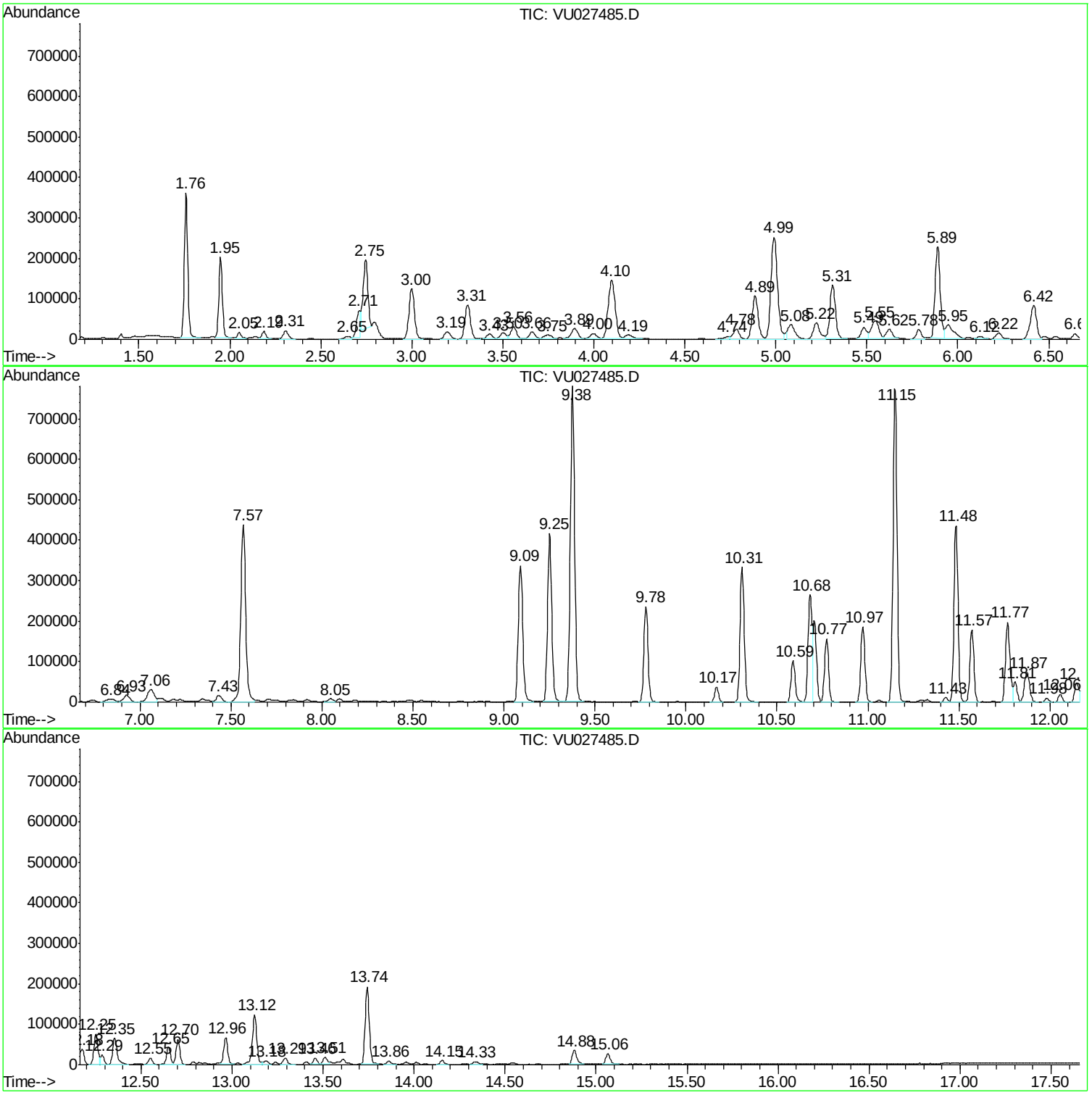
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU100618\
Data File : VU027485.D
Acq On : 06 Oct 2018 03:41
Operator : MD/SY
Sample : J5268-04 10X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-11

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.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\VU100618\
Data File : VU027485.D
Acq On : 06 Oct 2018 03:41
Operator : MD/SY
Sample : J5268-04 10X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-11

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

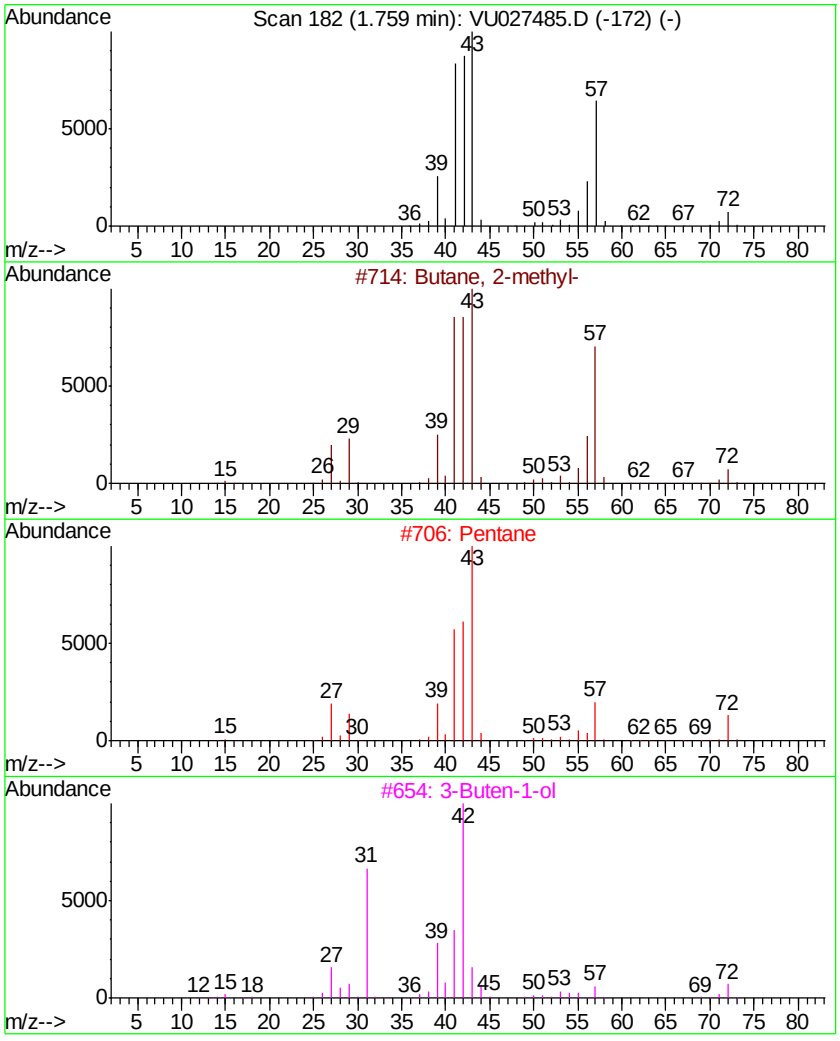
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

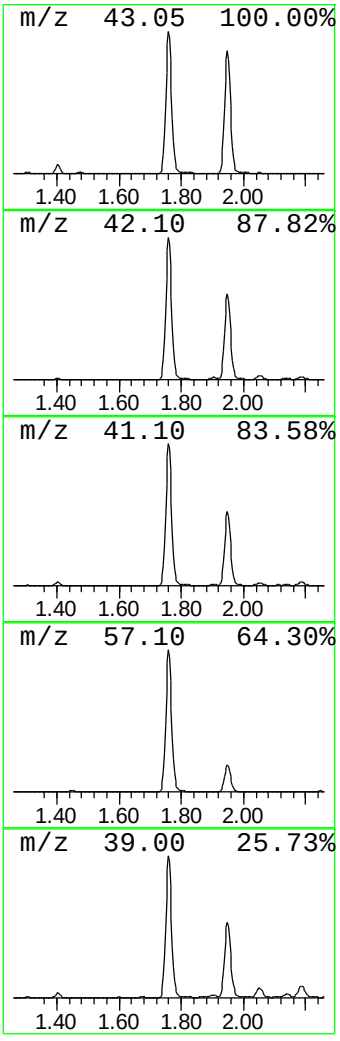
****

Peak Number 1 Butane, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.76	37.37 ug/l	473647	Pentafluorobenzene	4.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	91
2	Pentane	72	C5H12	000109-66-0	33
3	3-Buten-1-ol	72	C4H8O	000627-27-0	25
4	1-Butene	56	C4H8	000106-98-9	9
5	1-Propene, 2-methyl-	56	C4H8	000115-11-7	9





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100618\
Data File : VU027485.D
Acq On : 06 Oct 2018 03:41
Operator : MD/SY
Sample : J5268-04 10X
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-11

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

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

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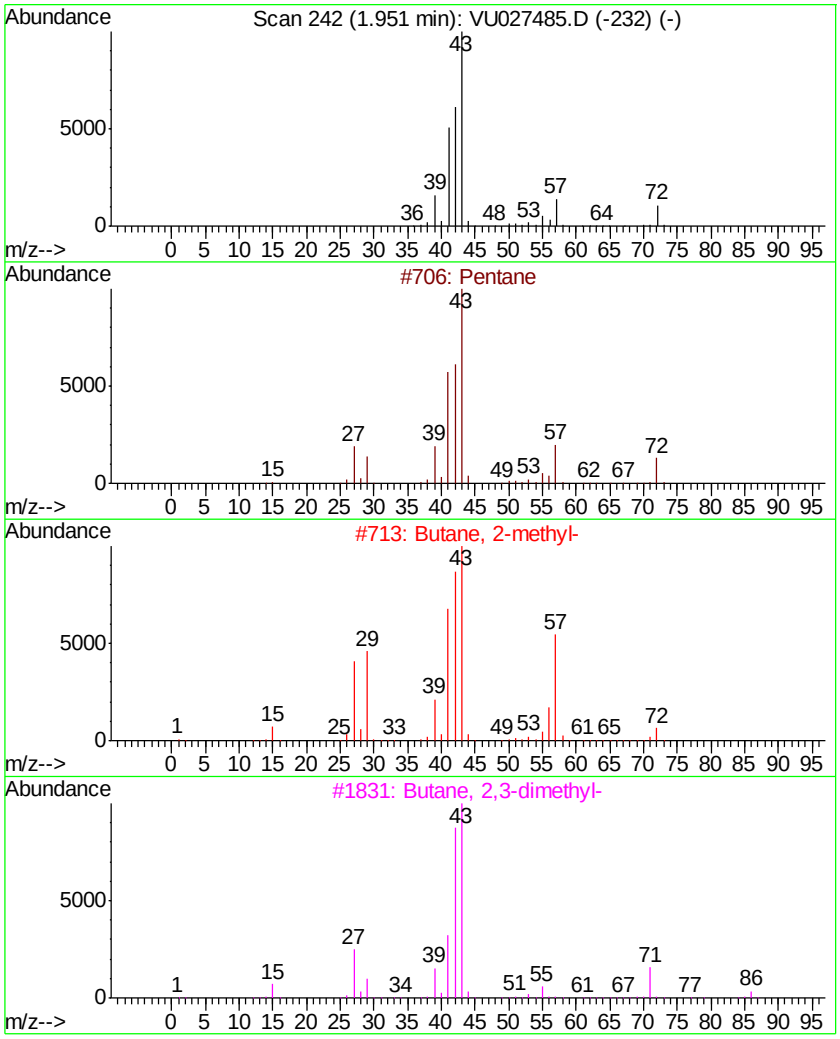
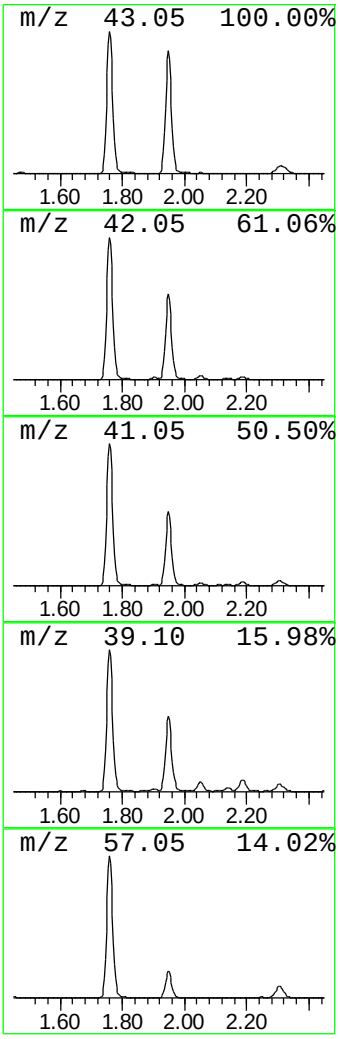
Peak Number 2 Pentane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.95	21.52 ug/l	272790	Pentafluorobenzene	4.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	91
2	Butane, 2-methyl-	72	C5H12	000078-78-4	45
3	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	36
4	Oxirane, ethyl-	72	C4H8O	000106-88-7	9
5	2-Hexanone, 5-methyl-	114	C7H14O	000110-12-3	9





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU100618\
Data File : VU027485.D
Acq On : 06 Oct 2018 03:41
Operator : MD/SY
Sample : J5268-04 10X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-11

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

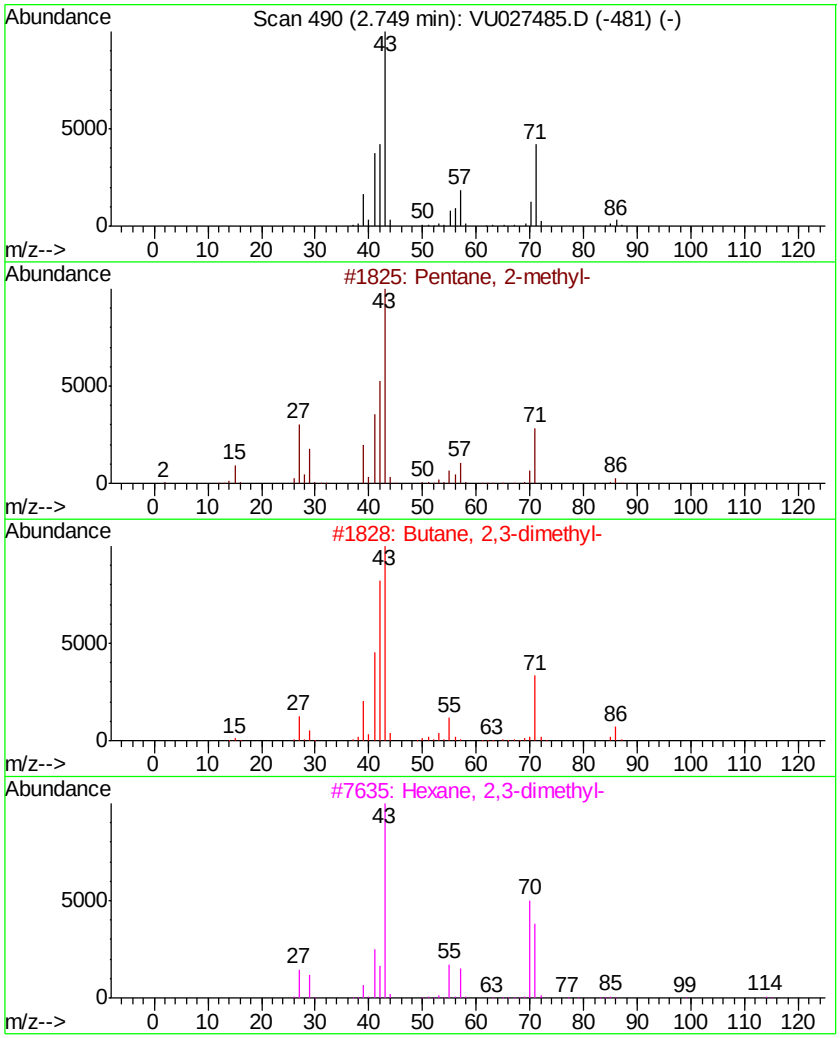
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

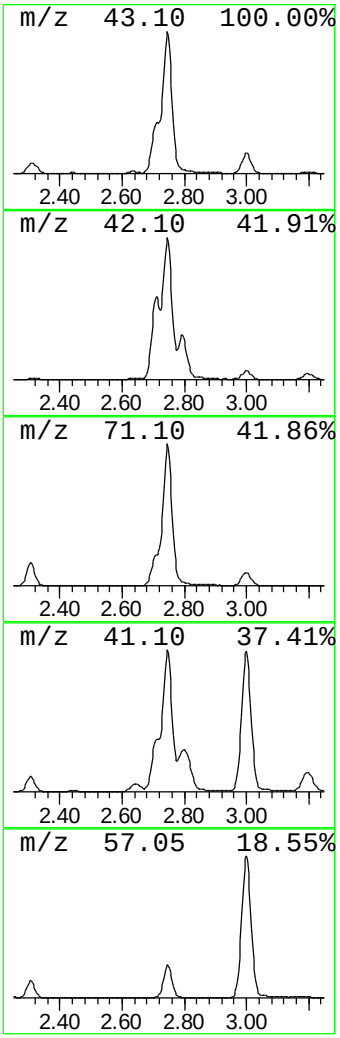
****

Peak Number 3 Pentane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.75	23.92 ug/l	303106	Pentafluorobenzene	4.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2-methyl-	86	C6H14	000107-83-5	64
2	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	47
3	Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	38
4	Pentane	72	C5H12	000109-66-0	38
5	1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	38





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU100618\
Data File : VU027485.D
Acq On : 06 Oct 2018 03:41
Operator : MD/SY
Sample : J5268-04 10X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-11

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

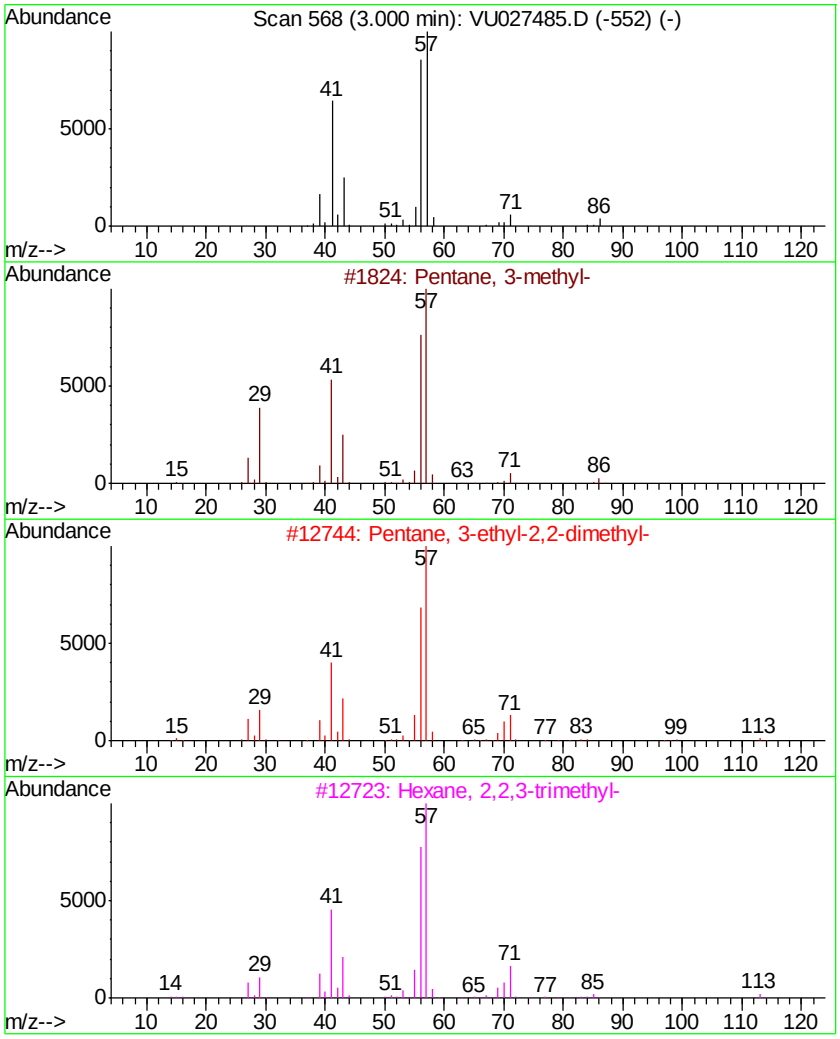
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

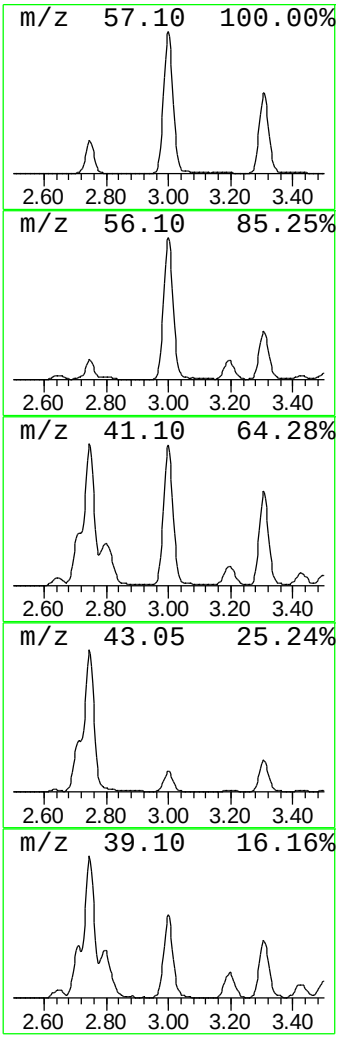
****

Peak Number 4 Pentane, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.00	20.79 ug/l	263478	Pentafluorobenzene	4.99

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64
3	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	64
4	Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	50
5	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	45





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU100618\
Data File : VU027485.D
Acq On : 06 Oct 2018 03:41
Operator : MD/SY
Sample : J5268-04 10X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-11

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

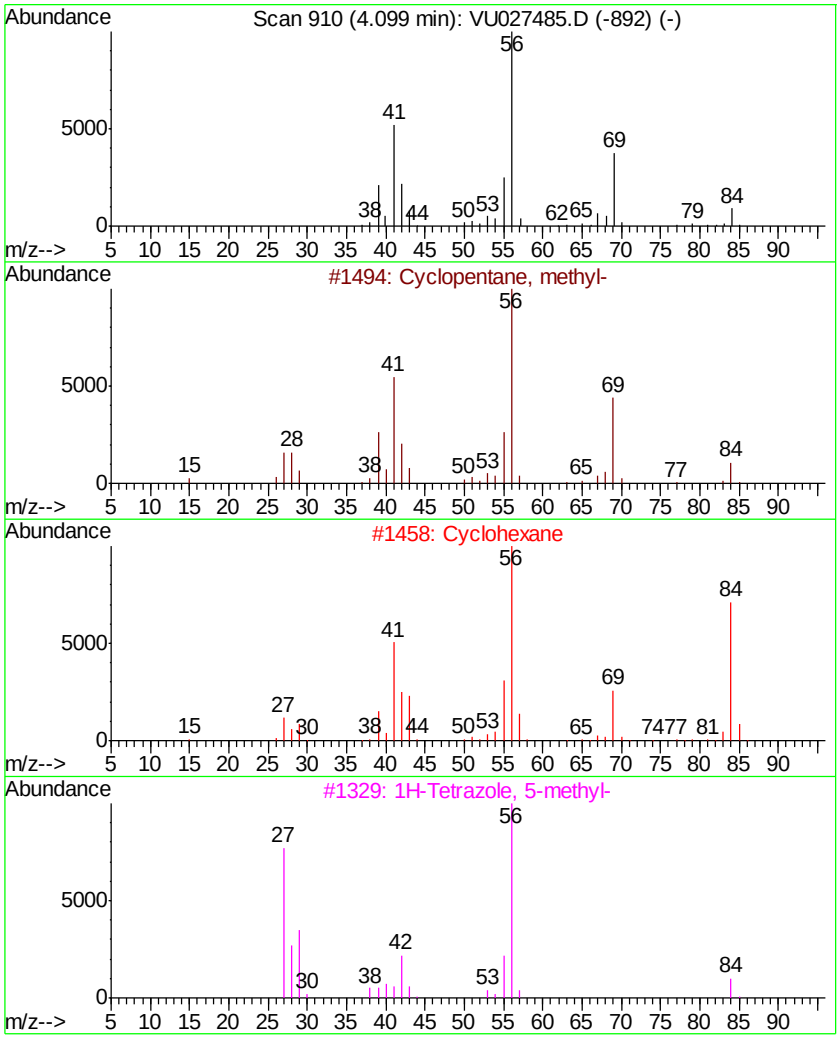
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

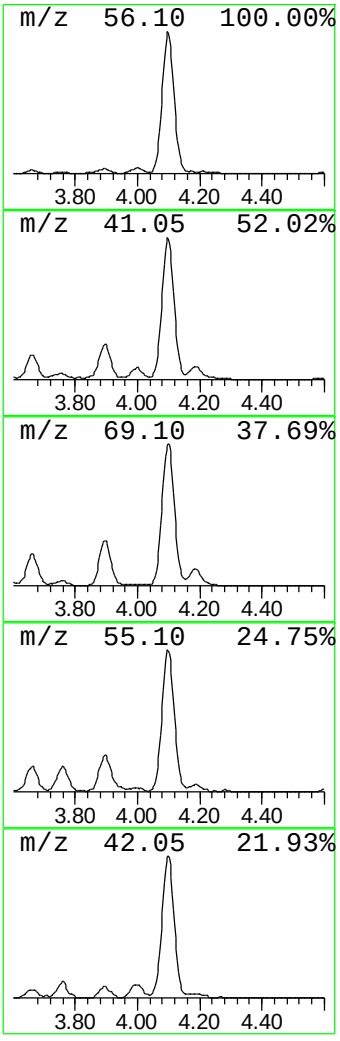
****

Peak Number 5 Cyclopentane, methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.10	30.61 ug/l	387917	Pentafluorobenzene	4.99

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	80
4	Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
5	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	50





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU100618\
Data File : VU027485.D
Acq On : 06 Oct 2018 03:41
Operator : MD/SY
Sample : J5268-04 10X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-11

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

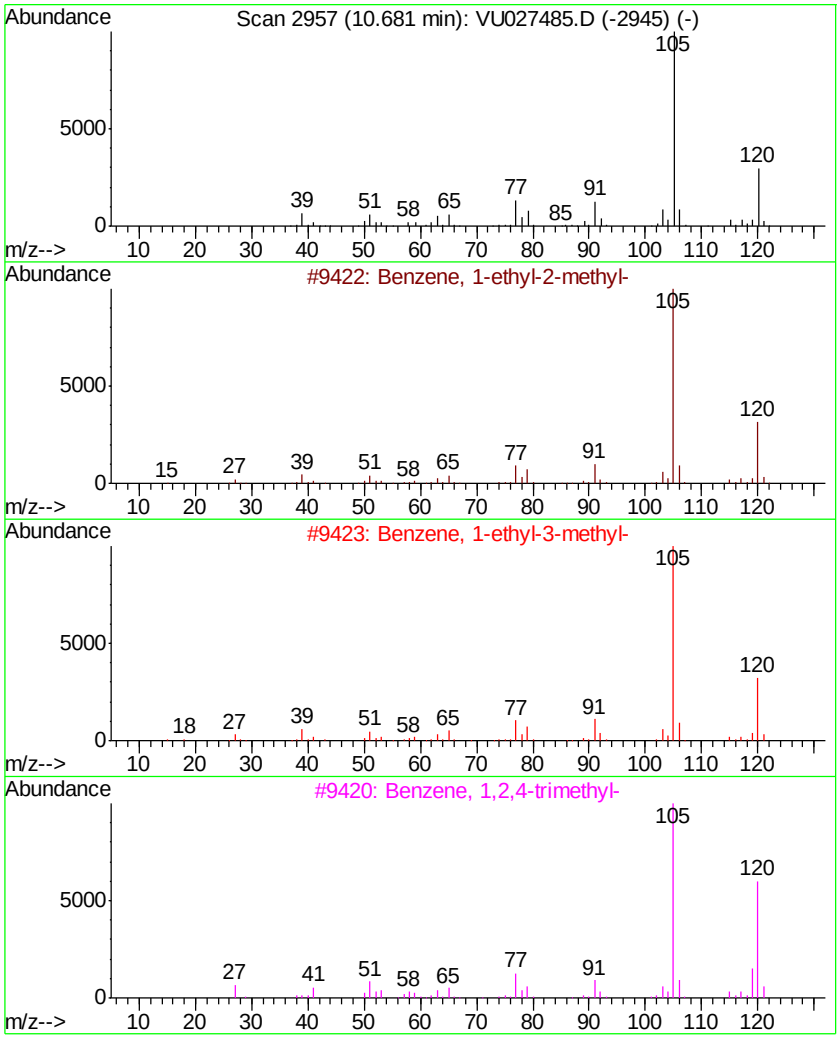
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

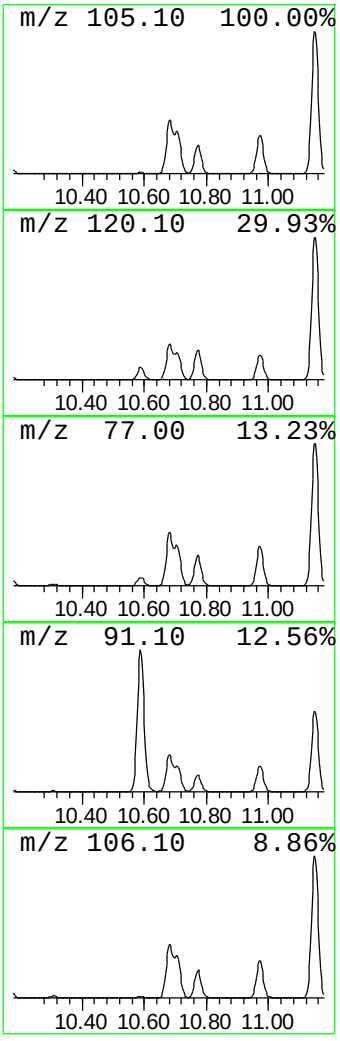
****

Peak Number 6 Benzene, 1-ethyl-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.68	31.20 ug/l	426429	1,4-Dichlorobenzene-d4	11.48

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-3-methyl-	120	C9H12	000620-14-4	95
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
5	Mesitylene	120	C9H12	000108-67-8	90





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU100618\
Data File : VU027485.D
Acq On : 06 Oct 2018 03:41
Operator : MD/SY
Sample : J5268-04 10X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-11

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

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

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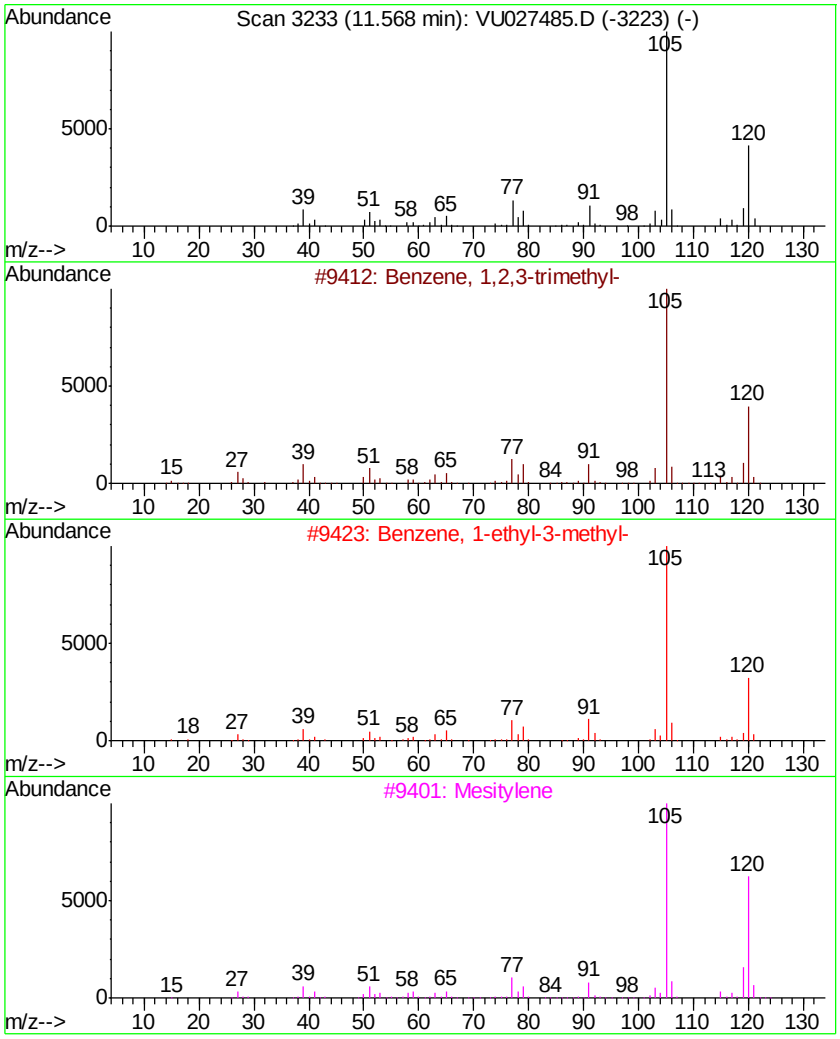
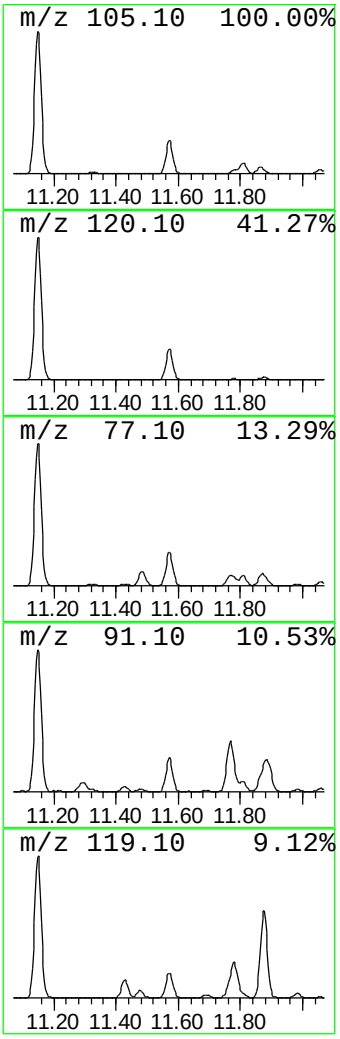
Peak Number 8 Benzene, 1,2,3-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	19.89 ug/l	271844	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU100618\
Data File : VU027485.D
Acq On : 06 Oct 2018 03:41
Operator : MD/SY
Sample : J5268-04 10X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-11

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

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

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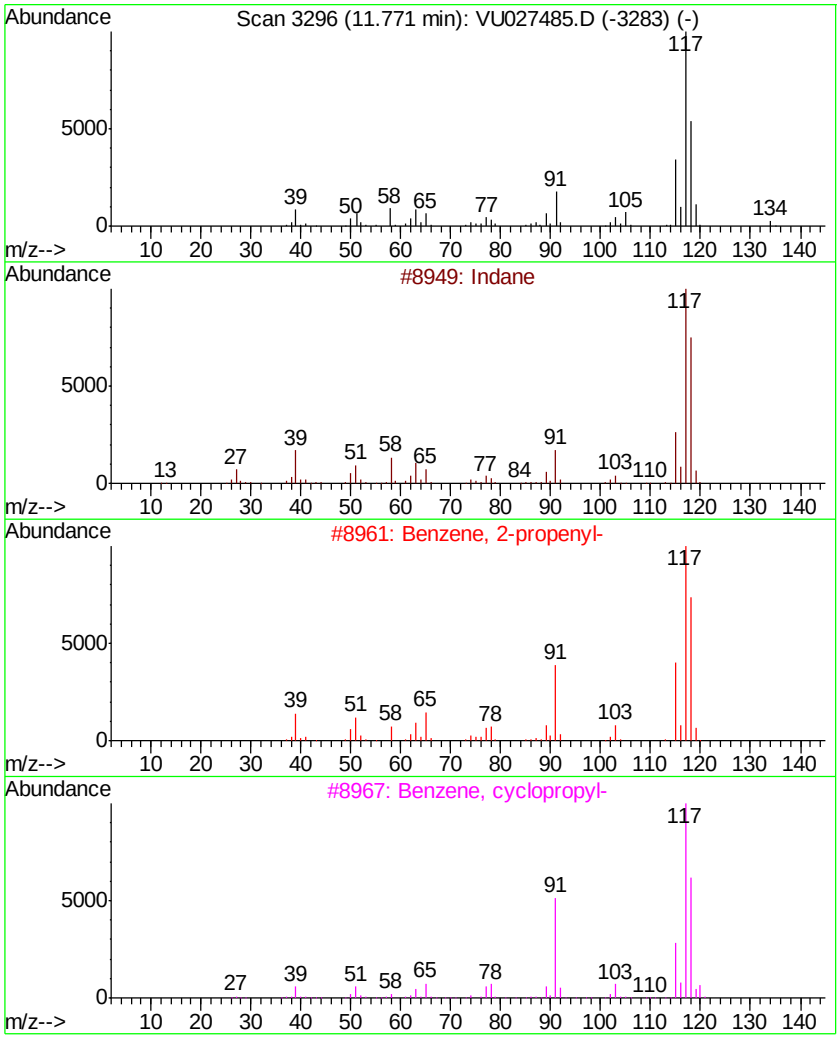
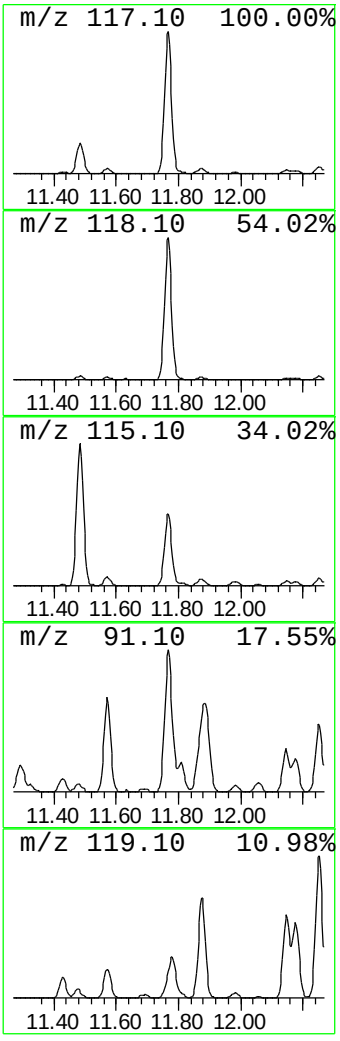
Peak Number 9 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.77	25.29 ug/l	345576	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	96
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	81
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	72
4	Deltacyclene	118	C9H10	007785-10-6	64
5	Benzene, 1-propenyl-	118	C9H10	000637-50-3	64





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU100618\
Data File : VU027485.D
Acq On : 06 Oct 2018 03:41
Operator : MD/SY
Sample : J5268-04 10X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-11

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

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

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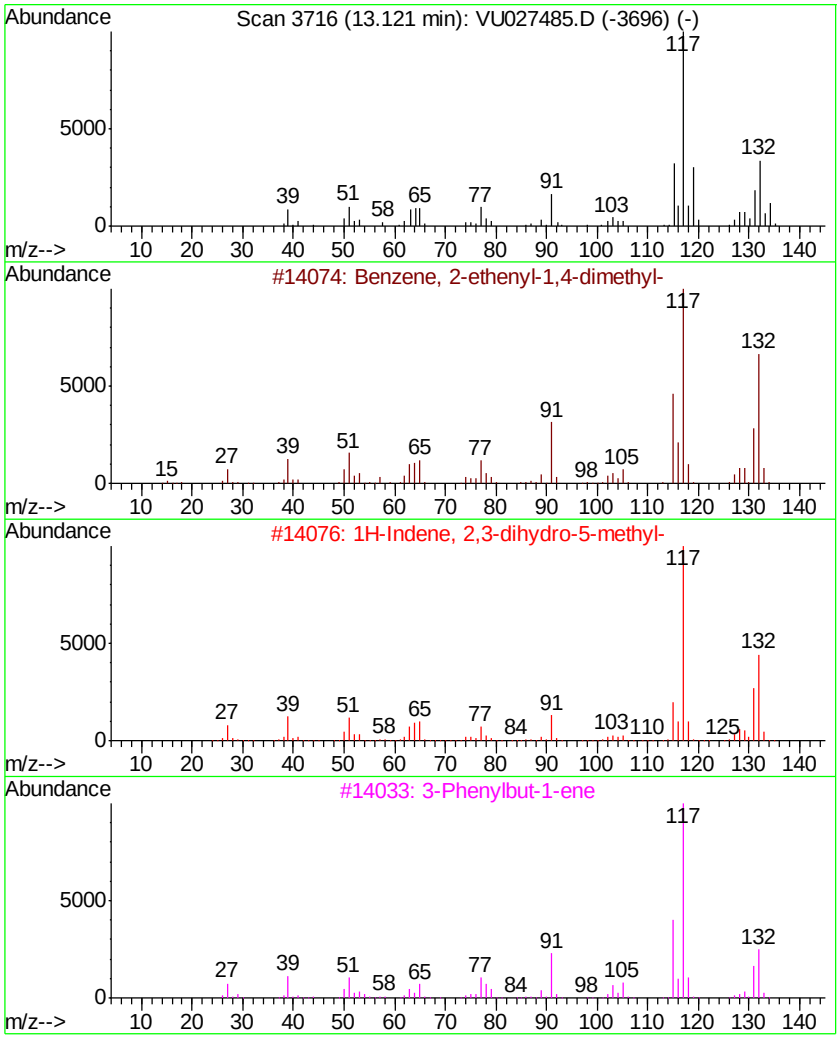
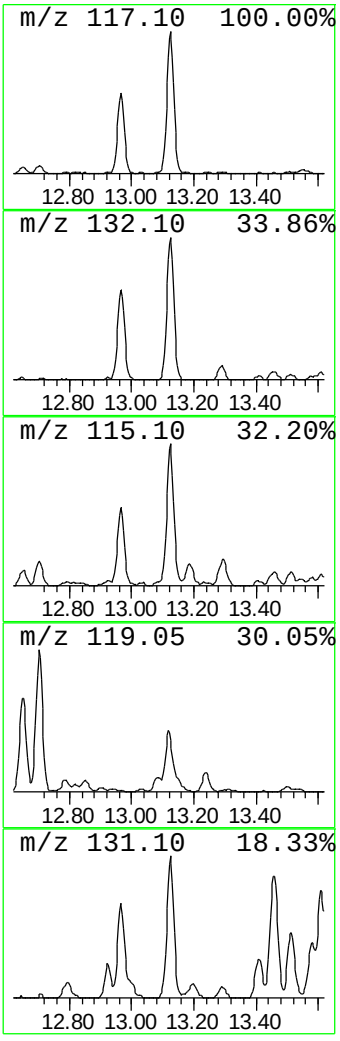
Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	15.65 ug/l	213870	1,4-Dichlorobenzene-d4	11.48

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	76
5	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	76





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_U\DATA\VU100618\
Data File : VU027485.D
Acq On : 06 Oct 2018 03:41
Operator : MD/SY
Sample : J5268-04 10X
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MW-11

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.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-methyl-	1.76	37.4	ug/l	473647	1	4.99	633673	50.0
Pentane	1.95	21.5	ug/l	272790	1	4.99	633673	50.0
Pentane, 2-methyl-	2.75	23.9	ug/l	303106	1	4.99	633673	50.0
Pentane, 3-methyl-	3.00	20.8	ug/l	263478	1	4.99	633673	50.0
Cyclopentane, met...	4.10	30.6	ug/l	387917	1	4.99	633673	50.0
Benzene, 1-ethyl-...	10.68	31.2	ug/l	426429	4	11.48	683311	50.0
Benzene, 1,2,3-tr...	11.57	19.9	ug/l	271844	4	11.48	683311	50.0
Indane	11.77	25.3	ug/l	345576	4	11.48	683311	50.0
Benzene, 2-etheny...	13.12	15.7	ug/l	213870	4	11.48	683311	50.0