

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU100618\
 Data File : VU027481.D
 Acq On : 06 Oct 2018 02:08
 Operator : MD/SY
 Sample : J5268-08
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 NW-4

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 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	195	rBV	89314	118606	12.11%	0.605%
2	1.948	232	241	253	rVB	47848	64427	6.58%	0.328%
3	2.321	343	357	374	rVB2	23250	48453	4.95%	0.247%
4	2.443	384	395	414	rBV	24668	44764	4.57%	0.228%
5	2.710	465	478	480	rBV2	38230	55880	5.71%	0.285%
6	2.746	482	489	501	rVB	101983	190124	19.42%	0.969%
7	3.000	553	568	592	rBV	93341	199330	20.36%	1.016%
8	3.189	612	627	645	rVB4	5492	13286	1.36%	0.068%
9	3.305	649	663	682	rBV2	41978	90333	9.23%	0.461%
10	4.000	861	879	892	rBV4	10268	27862	2.85%	0.142%
11	4.096	892	909	929	rVV2	95471	253028	25.84%	1.290%
12	4.276	955	965	982	rVB	8307	20446	2.09%	0.104%
13	4.736	1086	1108	1126	rBV6	8127	26404	2.70%	0.135%
14	4.887	1137	1155	1171	rBV	105953	245372	25.06%	1.251%
15	4.990	1171	1187	1203	rVV4	196842	484563	49.49%	2.470%
16	5.083	1206	1216	1238	rVB3	44427	122284	12.49%	0.623%
17	5.221	1244	1259	1276	rBV4	38975	118621	12.11%	0.605%
18	5.311	1276	1287	1315	rVV	131351	285590	29.17%	1.456%
19	5.482	1319	1340	1350	rVV3	60837	140721	14.37%	0.717%
20	5.553	1351	1362	1374	rVV5	67017	178366	18.22%	0.909%
21	5.627	1374	1385	1414	rVB2	52735	131911	13.47%	0.672%
22	5.787	1421	1435	1445	rBV4	12712	29096	2.97%	0.148%
23	5.887	1453	1466	1480	rBV	223206	453310	46.30%	2.311%
24	6.418	1610	1631	1658	rBV4	233939	599924	61.27%	3.058%
25	6.536	1659	1668	1677	rVV3	7076	12832	1.31%	0.065%
26	6.643	1691	1701	1717	rVV2	40237	74764	7.64%	0.381%
27	6.736	1721	1730	1742	rVB5	8750	17207	1.76%	0.088%
28	6.929	1772	1790	1808	rVB3	21422	53262	5.44%	0.272%
29	7.051	1810	1828	1843	rBV	41680	97838	9.99%	0.499%
30	7.446	1942	1951	1964	rVB6	9496	21436	2.19%	0.109%
31	7.565	1966	1988	2021	rBV	449565	881190	90.00%	4.492%
32	7.707	2021	2032	2040	rBV5	16506	35105	3.59%	0.179%
33	7.919	2087	2098	2115	rVB3	19681	41142	4.20%	0.210%
34	8.048	2118	2138	2148	rBV3	34344	71366	7.29%	0.364%

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35	8.179	2164	2179	2188	rBV4	7012	16983	1.73%	0.087%
36	8.491	2261	2276	2286	rBV7	17495	34795	3.55%	0.177%
37	8.546	2286	2293	2307	rVB4	10904	18663	1.91%	0.095%
38	9.089	2451	2462	2482	rVV	339688	580282	59.26%	2.958%
39	9.189	2483	2493	2502	rVV	20257	39464	4.03%	0.201%
40	9.250	2502	2512	2536	rVV	73510	125861	12.85%	0.642%
41	9.379	2536	2552	2567	rVB	71765	131267	13.41%	0.669%
42	9.781	2670	2677	2686	rVB2	9476	14373	1.47%	0.073%
43	10.167	2788	2797	2806	rBV2	20215	30587	3.12%	0.156%
44	10.308	2830	2841	2855	rBV	326672	516507	52.75%	2.633%
45	10.588	2918	2928	2947	rVB	132818	218357	22.30%	1.113%
46	10.707	2947	2965	2976	rVV	105608	226128	23.09%	1.153%
47	10.774	2976	2986	3002	rVB	96166	149159	15.23%	0.760%
48	10.970	3035	3047	3062	rBV	82631	132437	13.53%	0.675%
49	11.150	3091	3103	3117	rVB	292103	444153	45.36%	2.264%
50	11.292	3134	3147	3152	rBV	69183	106968	10.92%	0.545%
51	11.324	3152	3157	3173	rVB	67531	99827	10.20%	0.509%
52	11.427	3178	3189	3197	rBV	79714	120781	12.34%	0.616%
53	11.482	3197	3206	3225	rVV	457162	721903	73.73%	3.680%
54	11.572	3225	3234	3246	rVB2	23741	35736	3.65%	0.182%
55	11.687	3261	3270	3280	rVB2	11796	17263	1.76%	0.088%
56	11.784	3280	3300	3302	rBV5	296692	445531	45.50%	2.271%
57	11.810	3302	3308	3317	rVB	534365	787863	80.46%	4.017%
58	11.874	3317	3328	3349	rVB5	465798	979151	100.00%	4.992%
59	11.980	3349	3361	3372	rVB2	51278	79431	8.11%	0.405%
60	12.057	3372	3385	3398	rVB	190103	289830	29.60%	1.478%
61	12.147	3398	3413	3417	rBV	387068	568540	58.06%	2.898%
62	12.176	3418	3422	3433	rVB	365626	488818	49.92%	2.492%
63	12.250	3433	3445	3453	rBV	610487	922145	94.18%	4.701%
64	12.356	3467	3478	3486	rBV2	177073	324632	33.15%	1.655%
65	12.491	3509	3520	3527	rVB5	21185	36581	3.74%	0.186%
66	12.549	3527	3538	3550	rVB3	97068	174368	17.81%	0.889%
67	12.649	3551	3569	3577	rBV	375553	583150	59.56%	2.973%
68	12.703	3577	3586	3602	rVB	396418	606233	61.91%	3.091%
69	12.787	3602	3612	3618	rBV2	112733	178456	18.23%	0.910%
70	12.819	3618	3622	3628	rVB	45551	52324	5.34%	0.267%
71	12.922	3644	3654	3660	rBV3	67406	114141	11.66%	0.582%

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72	12.964	3660	3667	3675	rVB	196239	267886	27.36%	1.366%
73	13.028	3682	3687	3694	rVB2	30104	36666	3.74%	0.187%
74	13.092	3694	3707	3709	rBV5	89108	137785	14.07%	0.702%
75	13.121	3709	3716	3726	rVB3	269534	422503	43.15%	2.154%
76	13.237	3745	3752	3763	rVB	71098	100168	10.23%	0.511%
77	13.292	3763	3769	3792	rVB5	26399	54582	5.57%	0.278%
78	13.408	3792	3805	3812	rBV2	135988	220350	22.50%	1.123%
79	13.456	3812	3820	3828	rBV	211583	322257	32.91%	1.643%
80	13.510	3828	3837	3844	rBV3	307127	486329	49.67%	2.479%
81	13.578	3854	3858	3862	rBV	30157	29522	3.02%	0.151%
82	13.610	3862	3868	3874	rVV	116125	137699	14.06%	0.702%
83	13.742	3897	3909	3917	rBV	77282	125659	12.83%	0.641%
84	13.787	3917	3923	3937	rVV4	27649	54106	5.53%	0.276%
85	13.855	3937	3944	3958	rVB3	25114	42759	4.37%	0.218%
86	13.954	3959	3975	3986	rBV3	89660	177619	18.14%	0.905%
87	14.012	3986	3993	4004	rVV2	91993	142318	14.53%	0.726%
88	14.063	4004	4009	4023	rVB2	11540	18161	1.85%	0.093%
89	14.154	4026	4037	4048	rBV	115615	177659	18.14%	0.906%
90	14.334	4083	4093	4107	rBV	96476	196489	20.07%	1.002%
91	14.475	4129	4137	4148	rBV	49285	74401	7.60%	0.379%
92	14.539	4148	4157	4170	rVB2	63409	110945	11.33%	0.566%
93	14.607	4170	4178	4191	rVB3	13130	22797	2.33%	0.116%
94	14.687	4191	4203	4212	rBV5	16095	35279	3.60%	0.180%
95	14.877	4249	4262	4277	rBV	189168	313829	32.05%	1.600%
96	14.948	4277	4284	4294	rVB4	18874	31954	3.26%	0.163%
97	15.060	4310	4319	4333	rVV	90997	147466	15.06%	0.752%
98	15.916	4577	4585	4595	rBV3	13621	21697	2.22%	0.111%
99	16.060	4620	4630	4637	rBV2	19510	31202	3.19%	0.159%
100	16.102	4637	4643	4653	rVB3	8485	14026	1.43%	0.072%

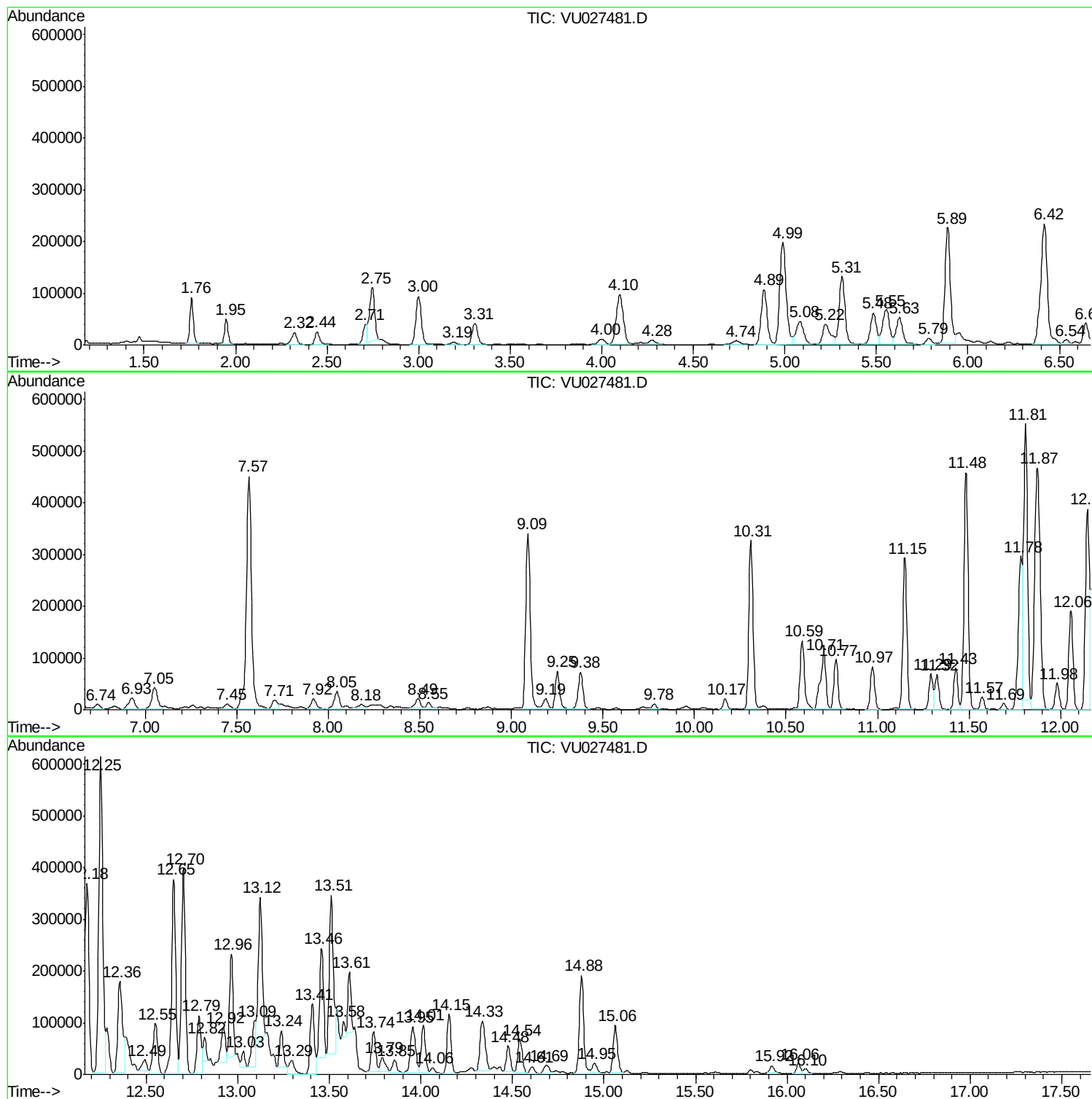
Sum of corrected areas: 19615644

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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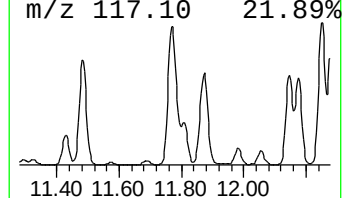
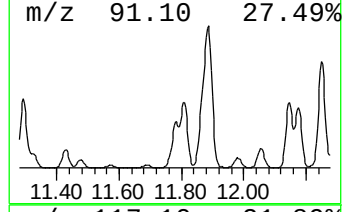
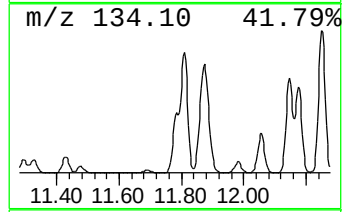
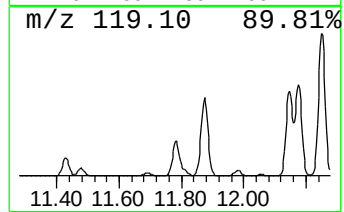
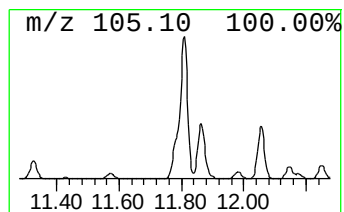
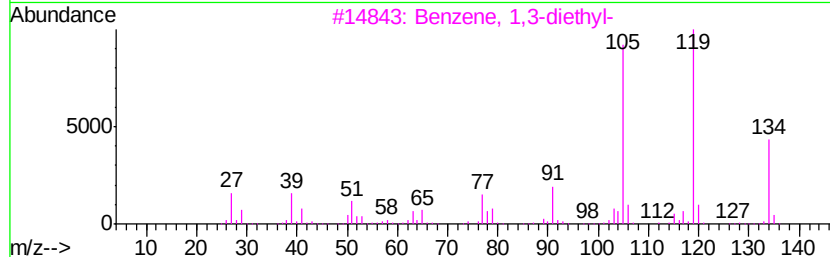
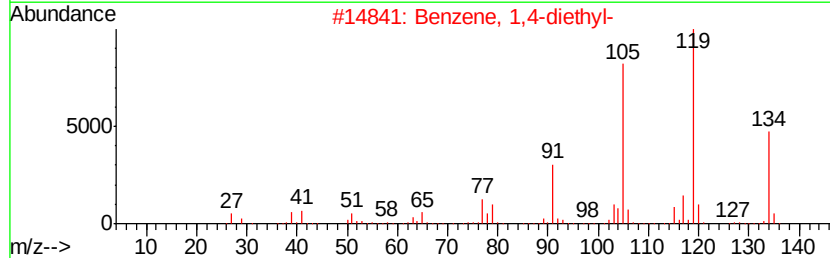
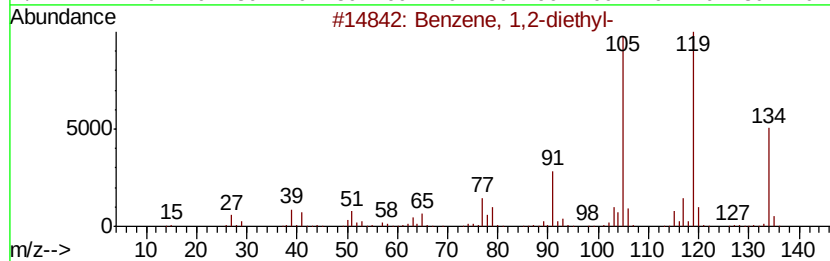
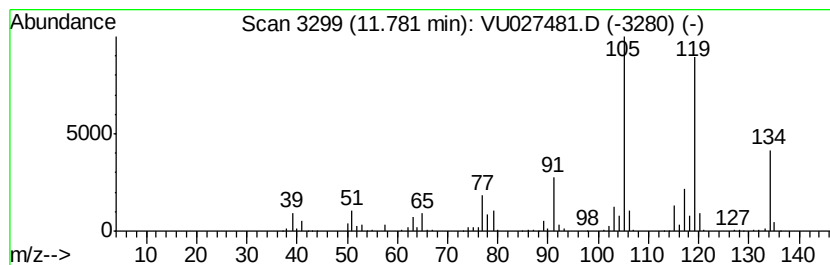
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1,2-diethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.78	30.86 ug/l	445531	1,4-Dichlorobenzene-d4	11.48	
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	95
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	93
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	87
5	1,3,8-p-Menthatriene	134	C10H14	018368-95-1	83



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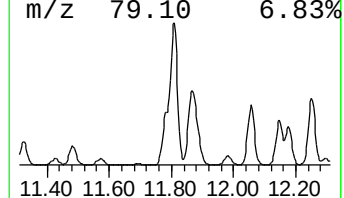
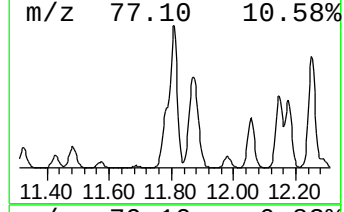
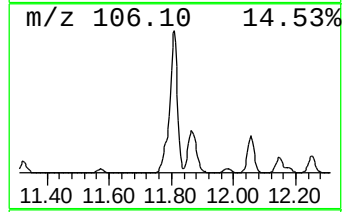
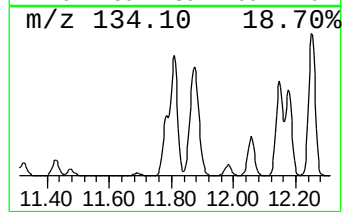
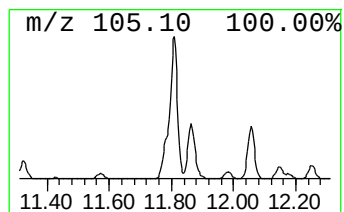
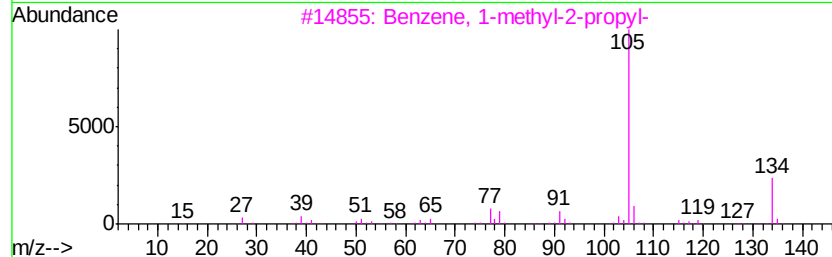
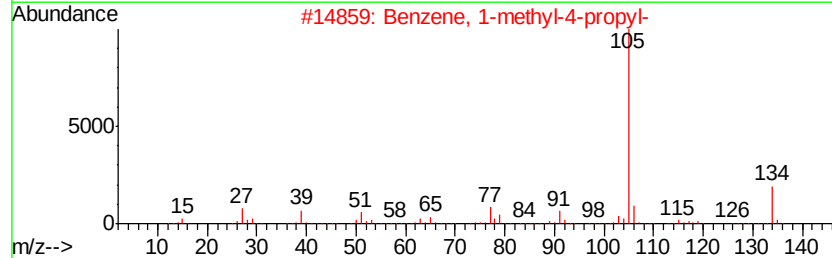
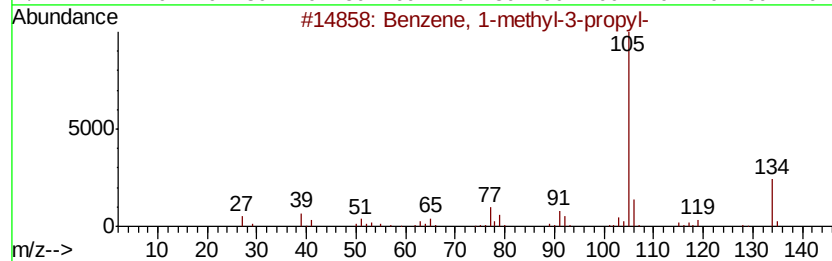
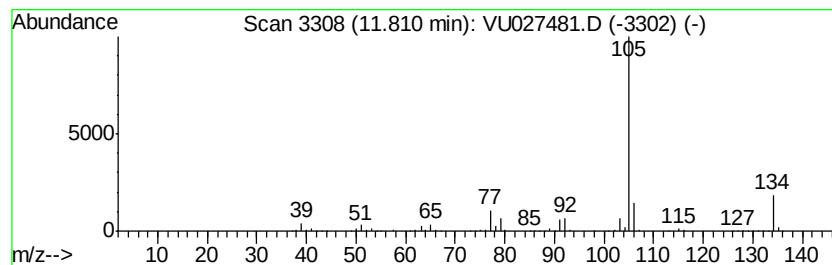
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TIC Library : C:\DATABASE\NIST11.L
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Peak Number 2 Benzene, 1-methyl-3-propyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.81	54.57 ug/l	787863	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	91
2		Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
3		Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	90
4		Benzeneacetaldehyde, 4-methyl-	134	C9H10O	000104-09-6	72
5		Benzeneacetaldehyde, .alpha.-met...	134	C9H10O	000093-53-8	72



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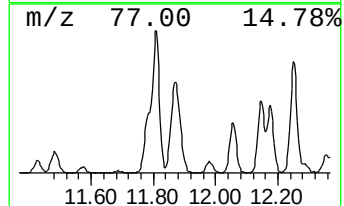
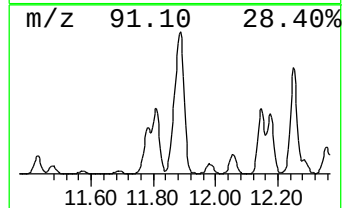
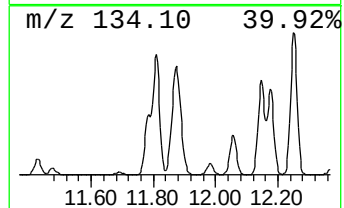
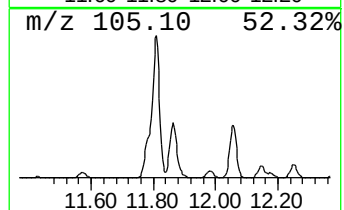
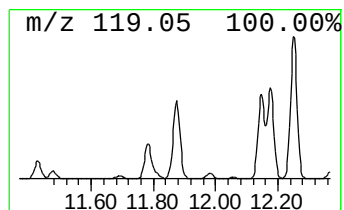
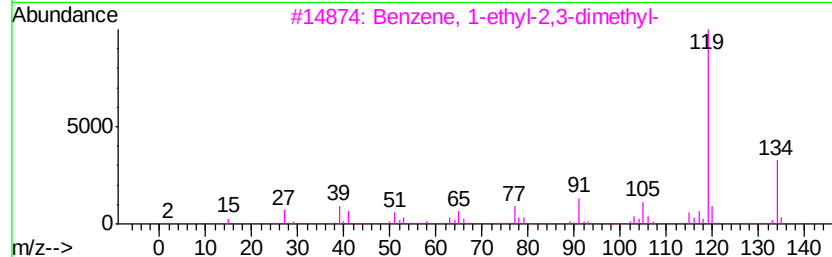
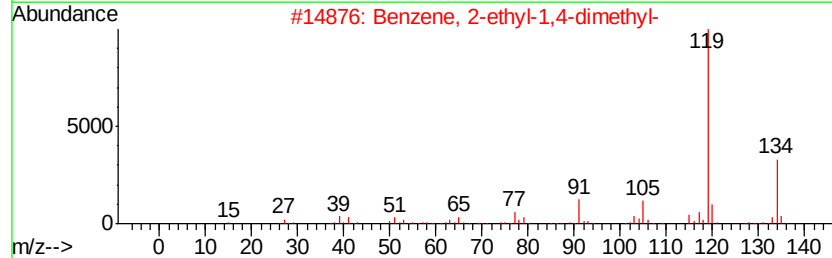
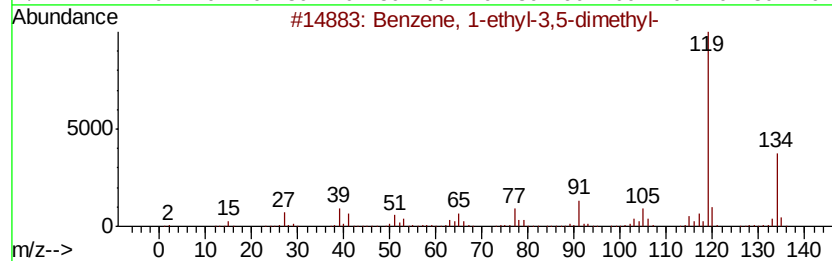
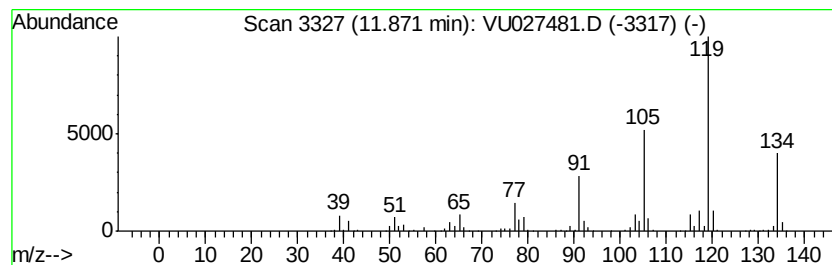
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TIC Integration Parameters: LSCINT.P

Peak Number 3 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.		
11.87	67.82 ug/l	979151	1,4-Dichlorobenzene-d4	11.48		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
2		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93
4		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	93
5		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100618\
Data File : VU027481.D
Acq On : 06 Oct 2018 02:08
Operator : MD/SY
Sample : J5268-08
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
NW-4

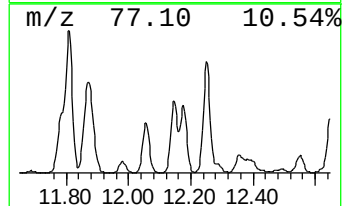
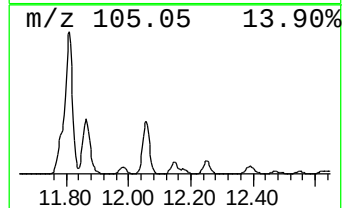
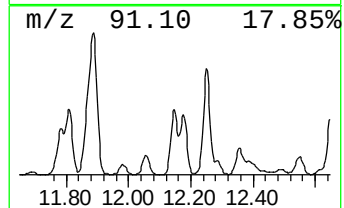
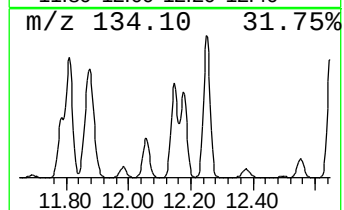
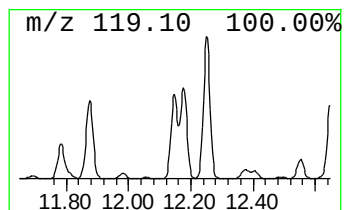
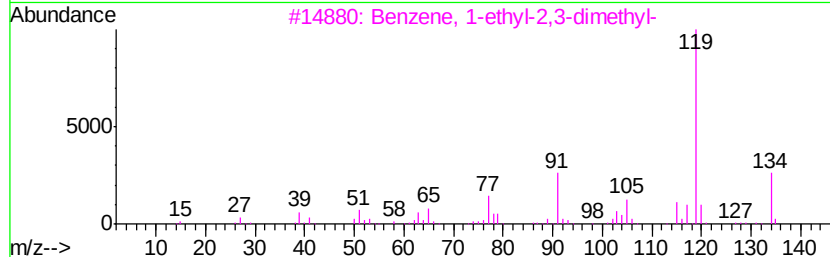
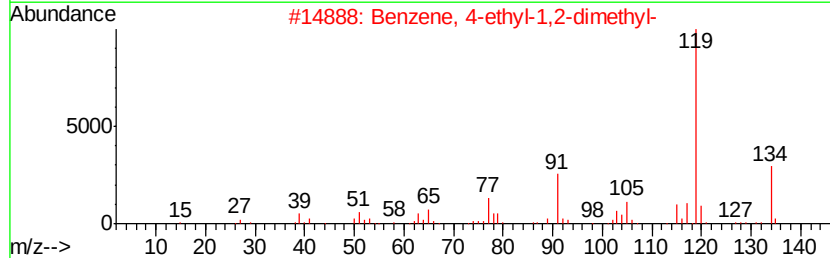
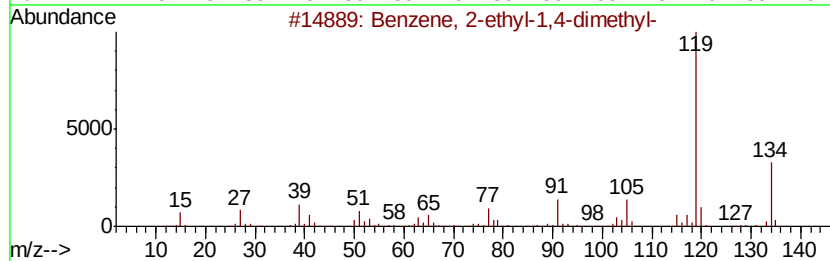
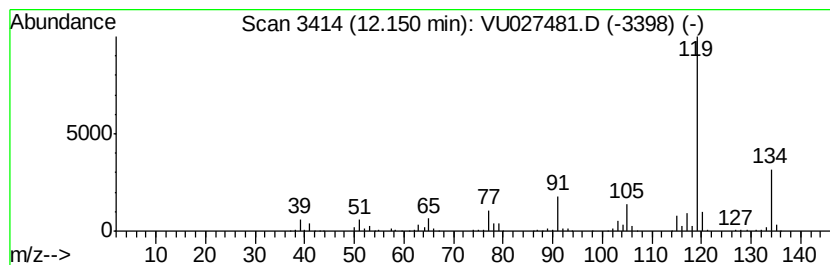
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.15	39.38 ug/l	568540	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
3		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
5		o-Cymene	134	C10H14	000527-84-4	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100618\
Data File : VU027481.D
Acq On : 06 Oct 2018 02:08
Operator : MD/SY
Sample : J5268-08
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
NW-4

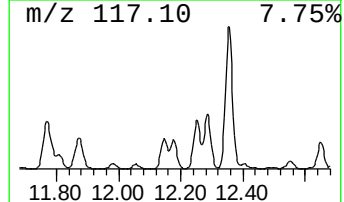
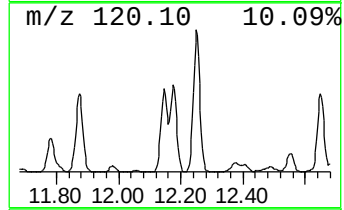
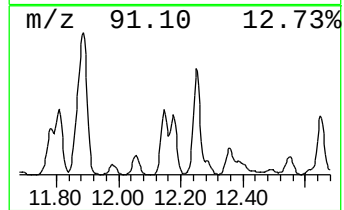
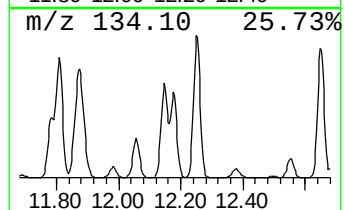
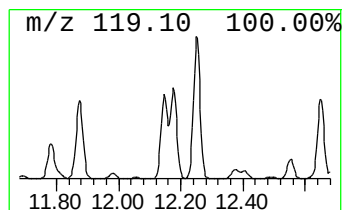
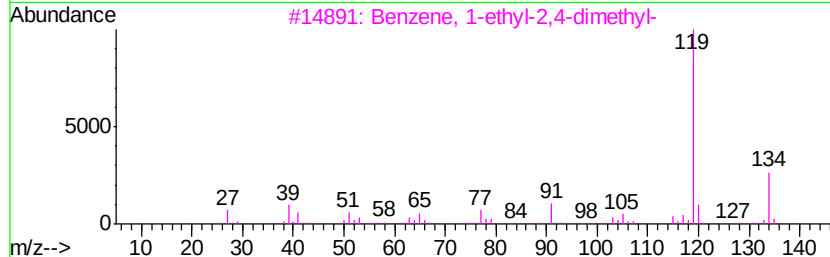
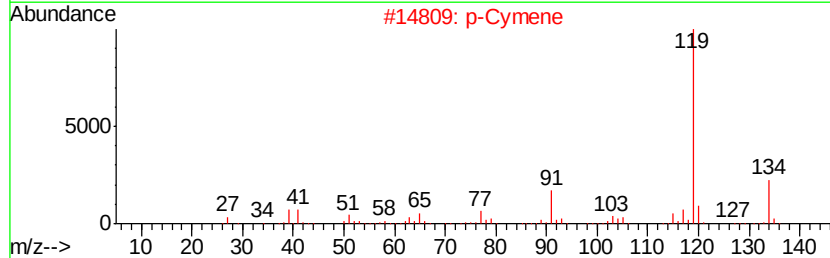
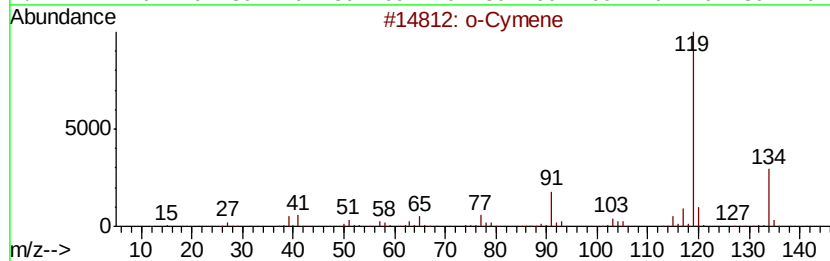
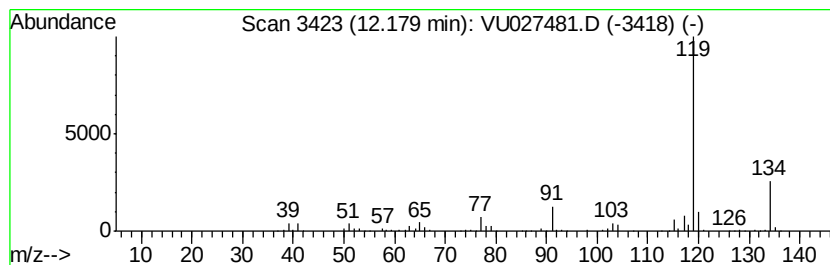
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Quant Title : SW846 8260

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

Peak Number 5 o-Cymene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.18	33.86 ug/l	488818	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	o-Cymene	134	C10H14	000527-84-4	94
2		p-Cymene	134	C10H14	000099-87-6	94
3		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	91
4		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	91
5		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100618\
Data File : VU027481.D
Acq On : 06 Oct 2018 02:08
Operator : MD/SY
Sample : J5268-08
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
NW-4

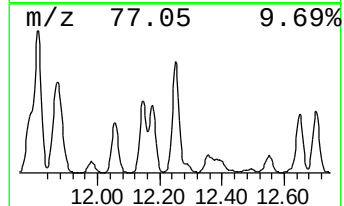
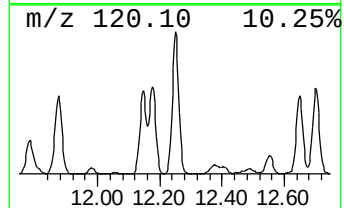
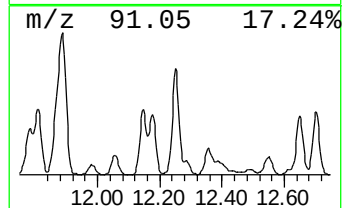
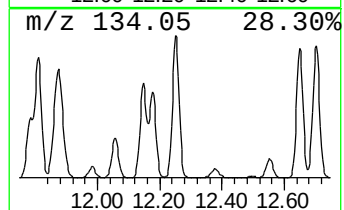
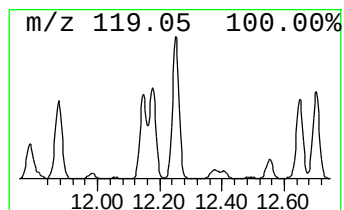
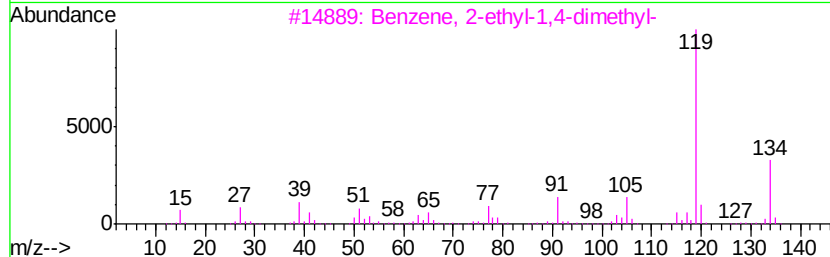
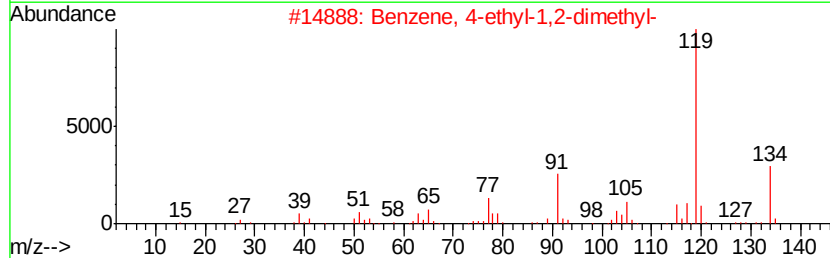
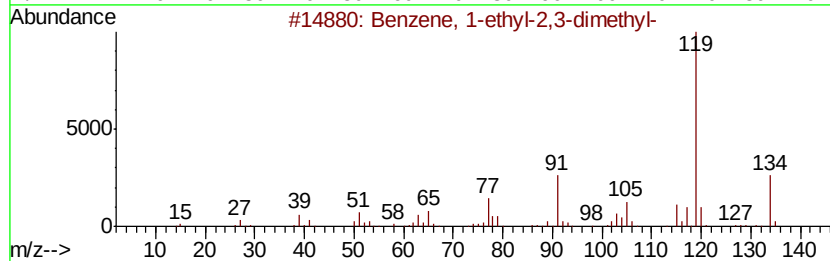
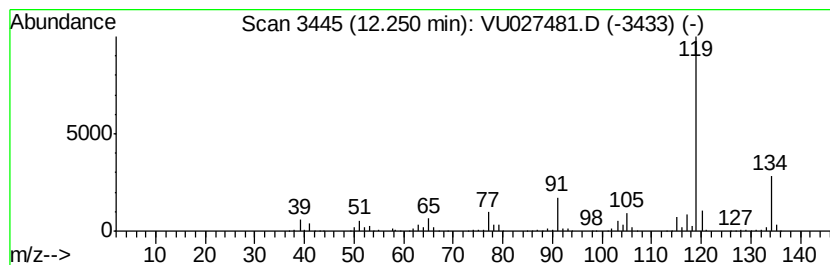
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.25	63.87 ug/l	922145	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
3		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5		o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100618\
Data File : VU027481.D
Acq On : 06 Oct 2018 02:08
Operator : MD/SY
Sample : J5268-08
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampled :
NW-4

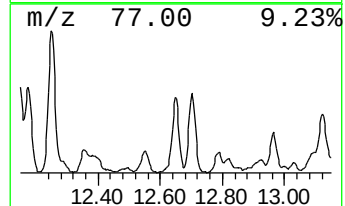
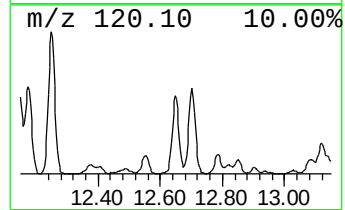
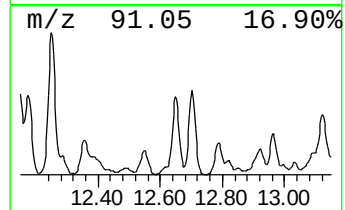
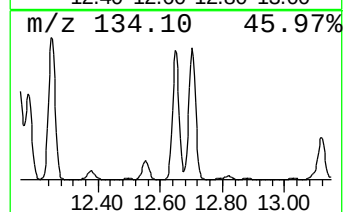
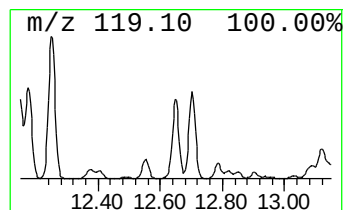
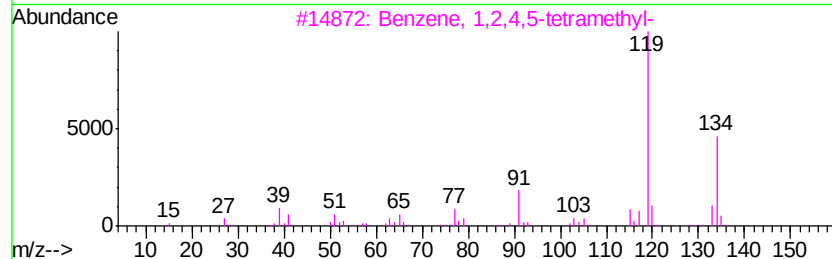
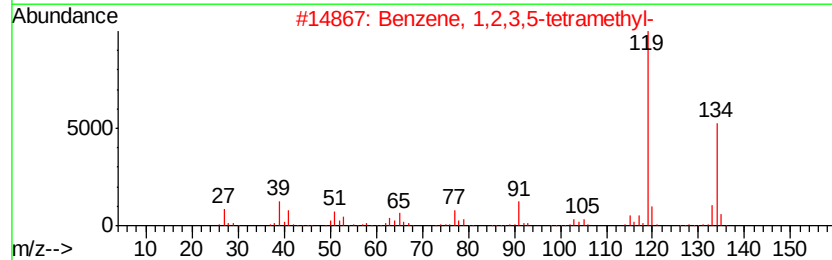
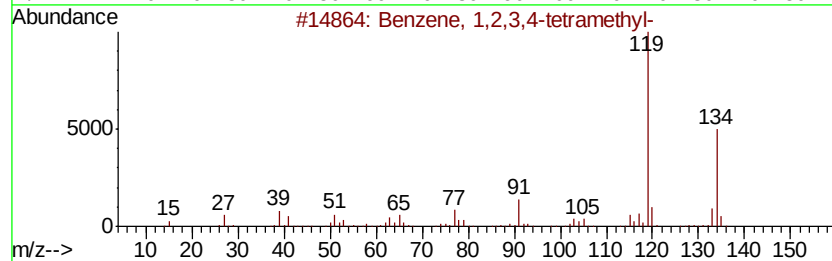
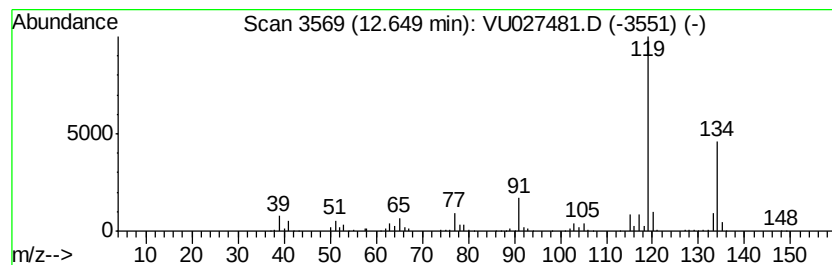
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.65	40.39 ug/l	583150	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
3		Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100618\
Data File : VU027481.D
Acq On : 06 Oct 2018 02:08
Operator : MD/SY
Sample : J5268-08
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
NW-4

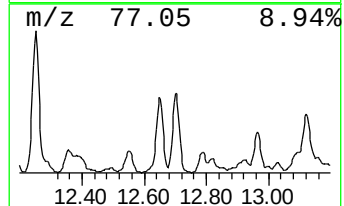
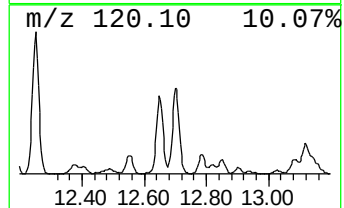
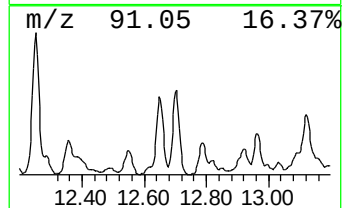
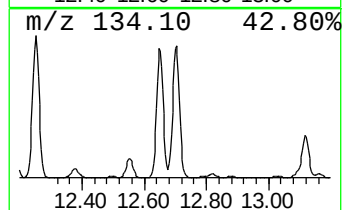
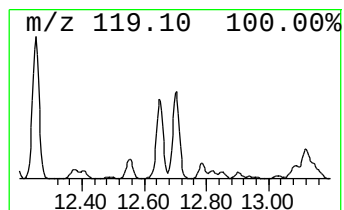
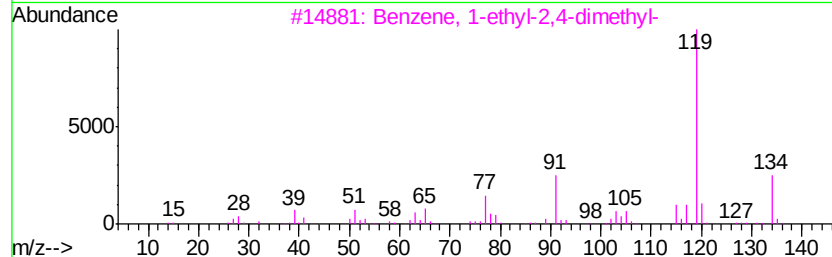
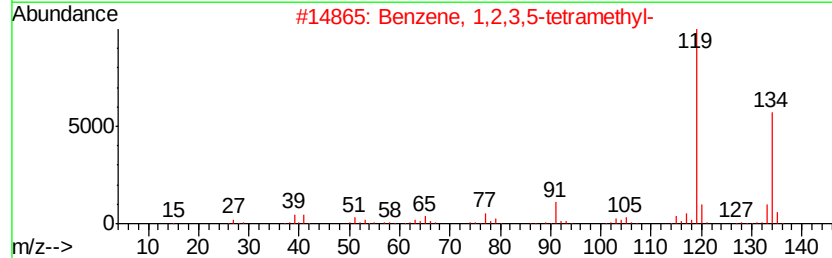
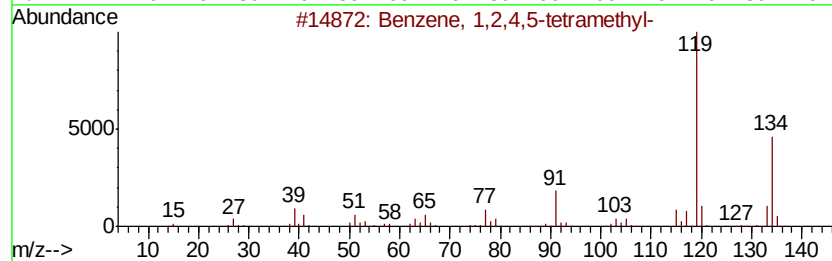
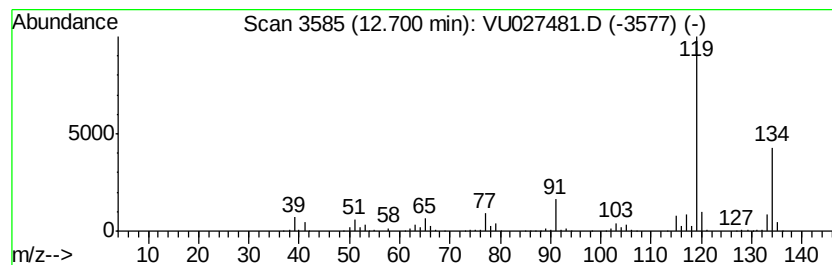
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

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

Peak Number 8 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.70	41.99 ug/l	606233	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
2		Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
3		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
4		Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
5		Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100618\
Data File : VU027481.D
Acq On : 06 Oct 2018 02:08
Operator : MD/SY
Sample : J5268-08
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
NW-4

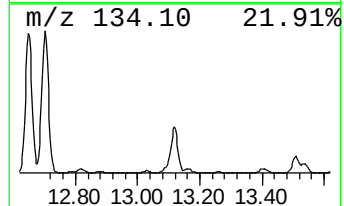
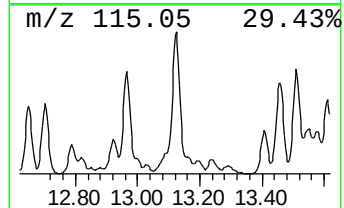
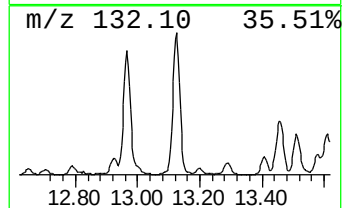
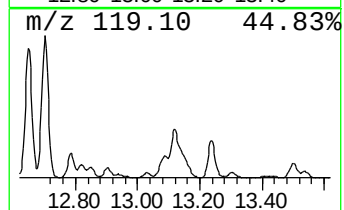
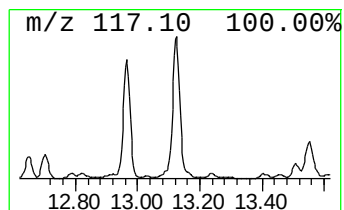
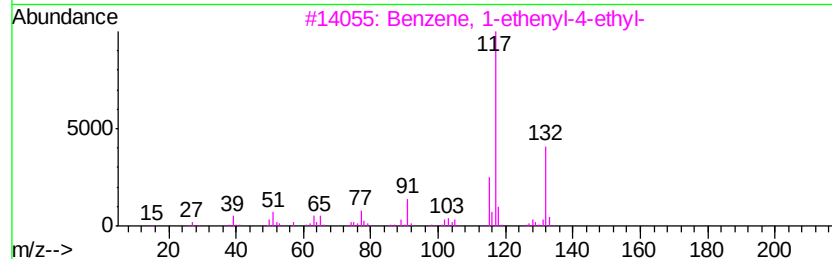
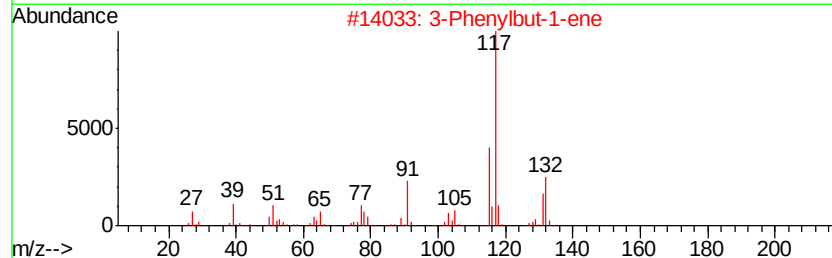
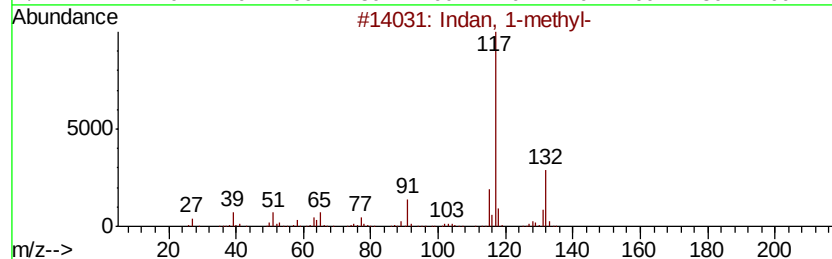
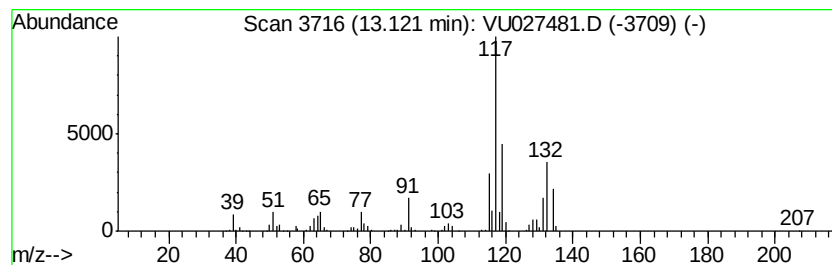
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

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

Peak Number 9 Indan, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.12	29.26 ug/l	422503	1,4-Dichlorobenzene-d4	11.48

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Indan, 1-methyl-	132	C10H12	000767-58-8	70
2		3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
3		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	64
4		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	64
5		Benzene, (2-methylcyclopropyl)-	132	C10H12	1000327-39-0	60



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA U\DATA\VU100618\
Data File : VU027481.D
Acq On : 06 Oct 2018 02:08
Operator : MD/SY
Sample : J5268-08
Misc : 5.0mL/MSVOA U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
NW-4

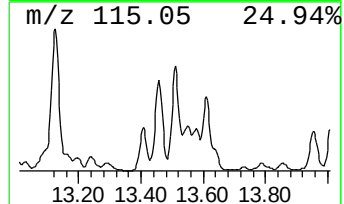
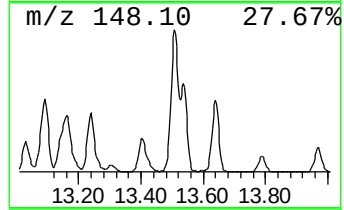
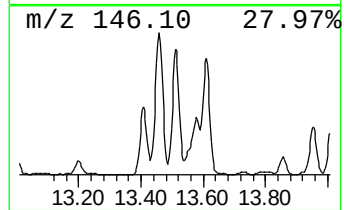
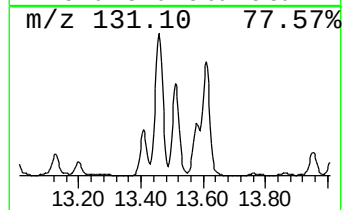
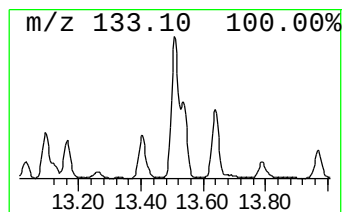
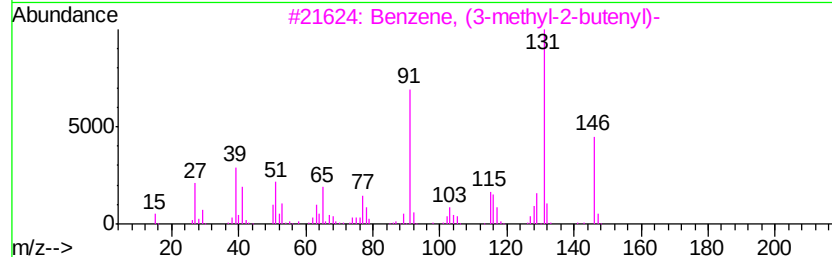
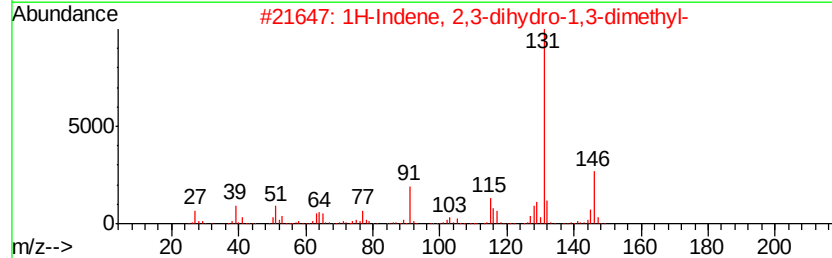
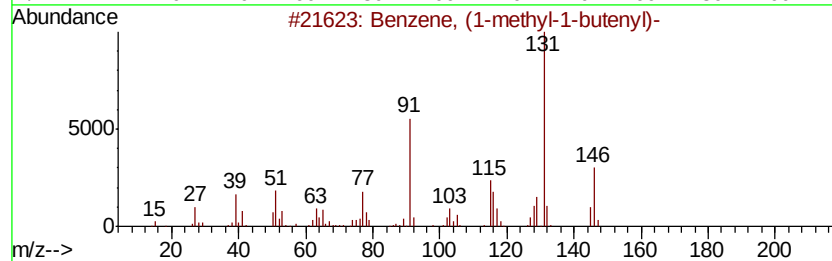
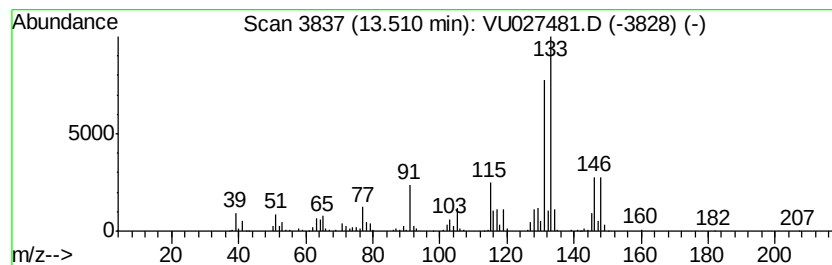
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

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

Peak Number 10 Benzene, (1-methyl-1-butenyl)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.51	33.68 ug/l	486329	1,4-Dichlorobenzene-d4	11.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	87
2		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	86
3		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	64
4		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	64
5		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_U\DATA\VU100618\
Data File : VU027481.D
Acq On : 06 Oct 2018 02:08
Operator : MD/SY
Sample : J5268-08
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
NW-4

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U100118W.M
Quant Title : SW846 8260

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1,2-diet...	11.78	30.9	ug/l	445531	4	11.48	721903	50.0
Benzene, 1-methyl...	11.81	54.6	ug/l	787863	4	11.48	721903	50.0
Benzene, 1-ethyl-...	11.87	67.8	ug/l	979151	4	11.48	721903	50.0
Benzene, 2-ethyl-...	12.15	39.4	ug/l	568540	4	11.48	721903	50.0
o-Cymene	12.18	33.9	ug/l	488818	4	11.48	721903	50.0
Benzene, 1-ethyl-...	12.25	63.9	ug/l	922145	4	11.48	721903	50.0
Benzene, 1,2,3,4-...	12.65	40.4	ug/l	583150	4	11.48	721903	50.0
Benzene, 1,2,4,5-...	12.70	42.0	ug/l	606233	4	11.48	721903	50.0
Indan, 1-methyl-	13.12	29.3	ug/l	422503	4	11.48	721903	50.0
Benzene, (1-methy...	13.51	33.7	ug/l	486329	4	11.48	721903	50.0