

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN092820\
 Data File : VN063718.D
 Acq On : 28 Sep 2020 20:20
 Operator : JC/MD
 Sample : L4178-06
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-6

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_N\METHODS\82N091620W.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	2.775	306	322	348	rBV	1045185	2127471	11.00%	0.927%
2	3.106	409	425	452	rVB3	261748	641716	3.32%	0.280%
3	3.531	539	557	576	rBV2	114841	281206	1.45%	0.123%
4	3.766	610	630	655	rVB4	84510	264365	1.37%	0.115%
5	4.547	834	873	932	rVB2	959710	4827321	24.95%	2.104%
6	5.007	987	1016	1052	rBV2	712837	2550929	13.19%	1.112%
7	5.331	1091	1117	1143	rBV4	60662	203248	1.05%	0.089%
8	5.881	1262	1288	1311	rBV2	201566	629919	3.26%	0.275%
9	6.023	1311	1332	1348	rVV4	144644	471751	2.44%	0.206%
10	6.328	1407	1427	1453	rBV6	225120	708073	3.66%	0.309%
11	6.476	1454	1473	1486	rBV4	141717	444021	2.30%	0.194%
12	6.582	1487	1506	1525	rBV	1610857	4958818	25.63%	2.162%
13	7.351	1732	1745	1770	rVB2	318166	865350	4.47%	0.377%
14	7.614	1783	1827	1844	rBV6	1773694	6090182	31.48%	2.655%
15	7.708	1844	1856	1878	rVB3	533363	1525625	7.89%	0.665%
16	7.839	1879	1897	1909	rBV	936359	2326010	12.02%	1.014%
17	7.997	1930	1946	1968	rBV3	356304	956385	4.94%	0.417%
18	8.126	1968	1986	1993	rBV5	804962	2046419	10.58%	0.892%
19	8.267	2020	2030	2052	rVB2	873269	2110873	10.91%	0.920%
20	8.376	2052	2064	2088	rVB5	88372	252271	1.30%	0.110%
21	8.550	2100	2118	2128	rBV3	1214297	2931499	15.15%	1.278%
22	8.714	2159	2169	2180	rVB	109034	202637	1.05%	0.088%
23	8.788	2180	2192	2197	rBV2	202847	389047	2.01%	0.170%
24	8.830	2198	2205	2241	rVB2	398781	897289	4.64%	0.391%
25	9.048	2243	2273	2286	rBV2	2577032	6896058	35.65%	3.006%
26	9.113	2286	2293	2307	rVB3	323110	631437	3.26%	0.275%
27	9.238	2321	2332	2344	rVB	487836	952836	4.93%	0.415%
28	9.334	2344	2362	2373	rBV3	248328	596574	3.08%	0.260%
29	9.402	2374	2383	2386	rBV4	165103	245494	1.27%	0.107%
30	9.486	2399	2409	2426	rVB3	870611	1768210	9.14%	0.771%
31	9.585	2428	2440	2446	rBV2	826200	1578265	8.16%	0.688%
32	9.630	2448	2454	2467	rVB5	718055	1406433	7.27%	0.613%
33	9.733	2479	2486	2500	rVB2	319727	656273	3.39%	0.286%
34	9.839	2500	2519	2539	rBV3	505978	1562965	8.08%	0.681%

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Integration Parameters: RTEINT.P

Integrator: RTE
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Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
 Title : SW846 8260

35	9.939	2539	2550	2559	rBV	623330	1186779	6.14%	0.517%
36	9.990	2560	2566	2575	rVB4	199288	342288	1.77%	0.149%
37	10.058	2575	2587	2603	rBV3	1667763	3750126	19.39%	1.635%
38	10.232	2632	2641	2643	rBV4	289349	416628	2.15%	0.182%
39	10.405	2686	2695	2709	rVB2	529440	946191	4.89%	0.412%
40	10.492	2710	2722	2739	rBV4	834979	1850169	9.56%	0.806%
41	10.691	2765	2784	2793	rBV2	246033	582342	3.01%	0.254%
42	10.765	2794	2807	2812	rVV3	314361	673357	3.48%	0.294%
43	10.801	2813	2818	2826	rVV4	387879	712354	3.68%	0.311%
44	10.849	2827	2833	2839	rVV4	289138	517669	2.68%	0.226%
45	10.891	2840	2846	2850	rVV	424082	675318	3.49%	0.294%
46	10.923	2851	2856	2865	rVV	606222	1049896	5.43%	0.458%
47	10.977	2865	2873	2884	rVV2	525271	971607	5.02%	0.424%
48	11.035	2884	2891	2899	rVV6	142405	258764	1.34%	0.113%
49	11.093	2900	2909	2917	rVB5	258309	474285	2.45%	0.207%
50	11.174	2918	2934	2955	rVB9	456144	1412894	7.30%	0.616%
51	11.273	2956	2965	2987	rBV	469942	1037571	5.36%	0.452%
52	11.383	2987	2999	3010	rBV2	1042810	1935052	10.00%	0.844%
53	11.486	3019	3031	3047	rVB	4564801	8039767	41.56%	3.505%
54	11.588	3048	3063	3069	rBV4	282329	736967	3.81%	0.321%
55	11.630	3071	3076	3085	rVB6	256064	405974	2.10%	0.177%
56	11.785	3117	3124	3132	rVV4	138529	244331	1.26%	0.107%
57	11.900	3148	3160	3174	rBV5	290368	895974	4.63%	0.391%
58	12.019	3191	3197	3205	rVB	199484	263704	1.36%	0.115%
59	12.093	3212	3220	3227	rBV4	352665	577878	2.99%	0.252%
60	12.138	3229	3234	3243	rVB6	254117	376947	1.95%	0.164%
61	12.225	3251	3261	3270	rBV	1445296	2313836	11.96%	1.009%
62	12.376	3300	3308	3317	rVB	718750	1089120	5.63%	0.475%
63	12.566	3352	3367	3378	rVB	3845158	6198714	32.04%	2.702%
64	12.707	3402	3411	3421	rVB	1423550	2386938	12.34%	1.040%
65	12.865	3448	3460	3475	rBV	3009530	5067193	26.19%	2.209%
66	12.936	3477	3482	3496	rVB4	323390	540658	2.79%	0.236%
67	13.016	3496	3507	3518	rBV	3251732	5137645	26.56%	2.240%
68	13.141	3533	3546	3556	rBV4	654662	1611575	8.33%	0.702%
69	13.264	3577	3584	3590	rBV2	315970	444732	2.30%	0.194%
70	13.318	3592	3601	3607	rBV	873268	1512702	7.82%	0.659%
71	13.485	3643	3653	3657	rBV	3364704	5055760	26.14%	2.204%

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72	13.518	3657	3663	3672	rVV	6042572	9947405	51.42%	4.336%
73	13.566	3672	3678	3692	rVB2	1892292	3127735	16.17%	1.363%
74	13.646	3693	3703	3712	rBV4	1079877	1797120	9.29%	0.783%
75	13.714	3718	3724	3735	rVB	346737	509673	2.63%	0.222%
76	13.778	3736	3744	3759	rVB	560852	877306	4.54%	0.382%
77	13.865	3759	3771	3781	rBV	8884839	14072525	72.75%	6.134%
78	13.919	3781	3788	3794	rVV	1262232	1910699	9.88%	0.833%
79	13.968	3794	3803	3815	rVB2	5821879	9591645	49.58%	4.181%
80	14.183	3853	3870	3878	rBV	5321288	8961019	46.32%	3.906%
81	14.222	3878	3882	3890	rVB	1395839	1779762	9.20%	0.776%
82	14.267	3890	3896	3903	rBV2	992894	1374922	7.11%	0.599%
83	14.450	3944	3953	3963	rBV2	4121021	6498666	33.59%	2.833%
84	14.498	3963	3968	3976	rVB2	845850	1173316	6.07%	0.511%
85	14.566	3976	3989	4000	rBV3	9954410	19344292	100.00%	8.432%
86	14.624	4000	4007	4018	rVB3	1189265	1957801	10.12%	0.853%
87	14.714	4028	4035	4043	rBV	540490	759090	3.92%	0.331%
88	14.823	4051	4069	4076	rBV4	2501266	5459736	28.22%	2.380%
89	14.929	4091	4102	4123	rVB2	2549369	5488133	28.37%	2.392%
90	15.103	4136	4156	4168	rBV	3810845	7153966	36.98%	3.118%
91	15.212	4180	4190	4197	rVV2	720384	1322242	6.84%	0.576%
92	15.260	4198	4205	4231	rVB	706570	1475052	7.63%	0.643%
93	15.386	4232	4244	4257	rBV	939315	1733749	8.96%	0.756%
94	15.543	4279	4293	4313	rBV2	592954	1581171	8.17%	0.689%
95	15.665	4322	4331	4340	rBV	228847	399059	2.06%	0.174%
96	15.736	4342	4353	4364	rVB4	357010	743556	3.84%	0.324%
97	15.874	4383	4396	4407	rBV3	180741	384370	1.99%	0.168%
98	16.122	4462	4473	4501	rVB	448263	1004925	5.19%	0.438%
99	16.312	4517	4532	4557	rVB	1054243	2287767	11.83%	0.997%

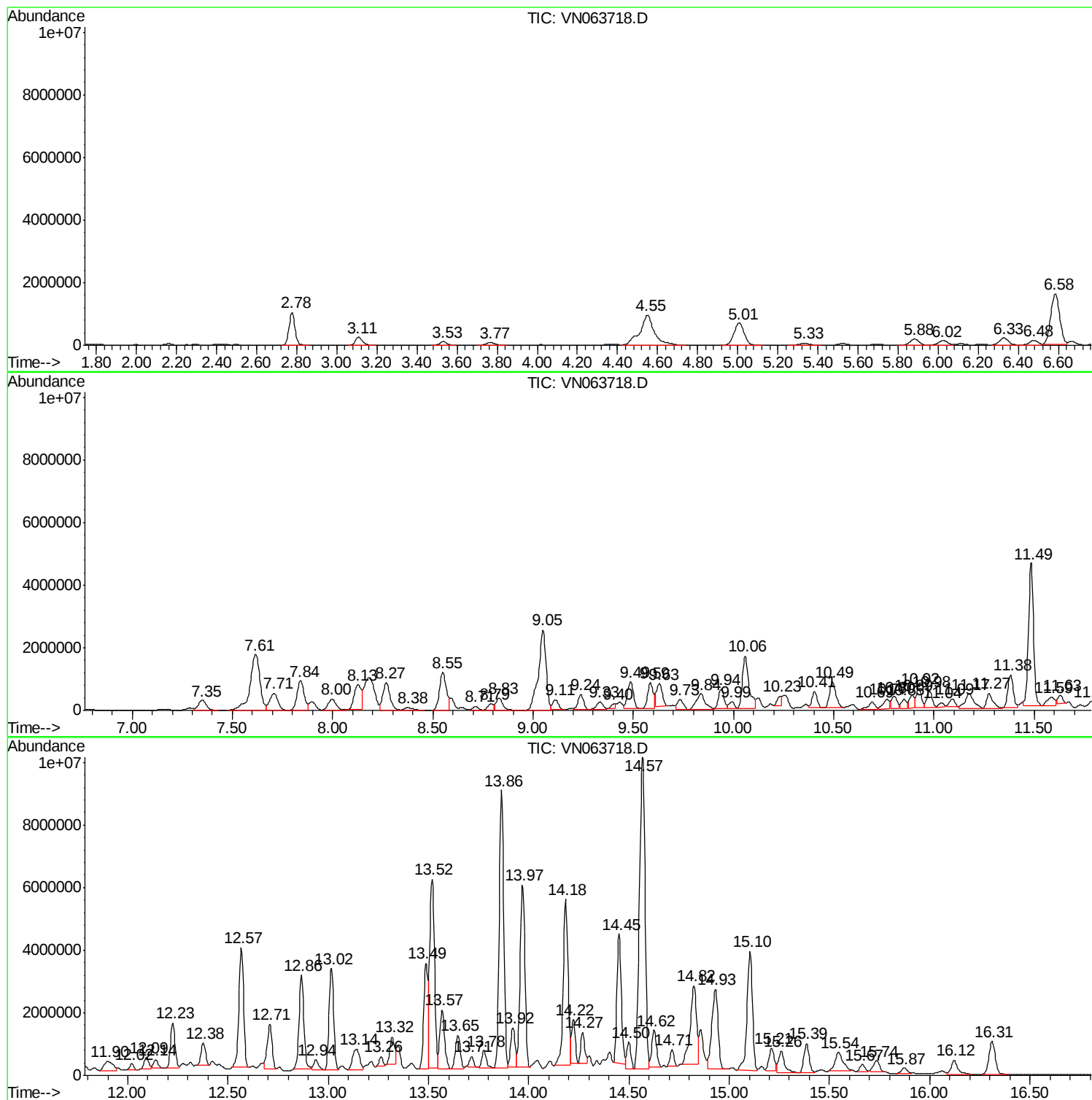
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Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
Quant Title : SW846 8260

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



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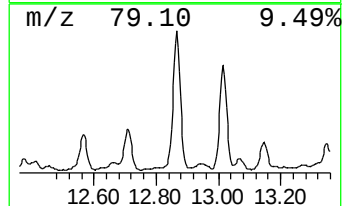
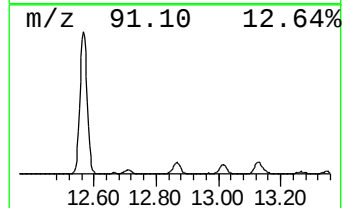
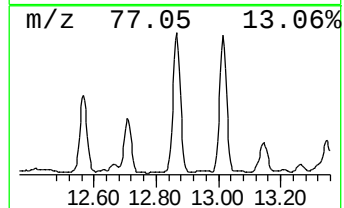
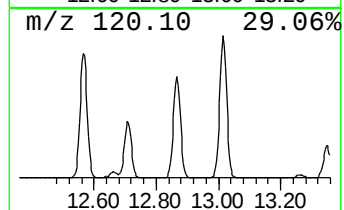
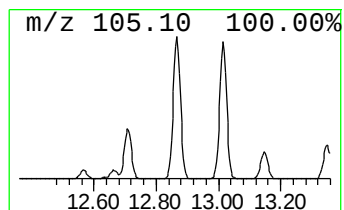
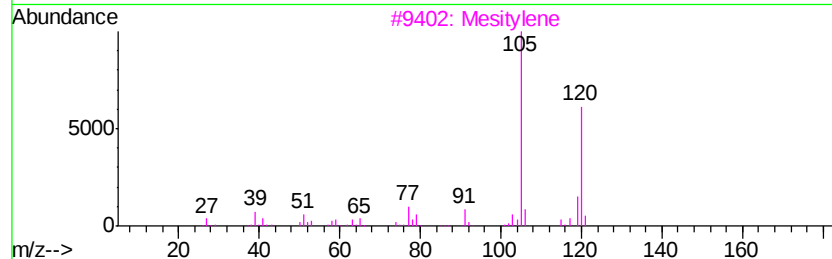
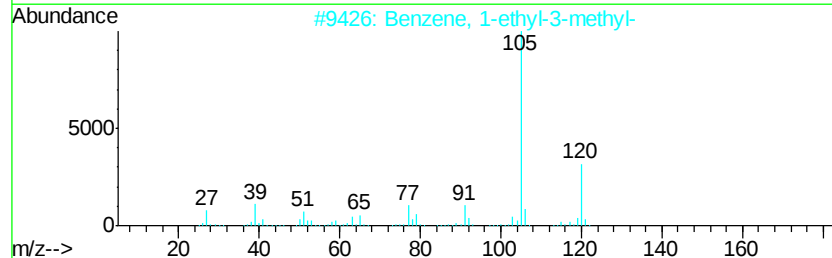
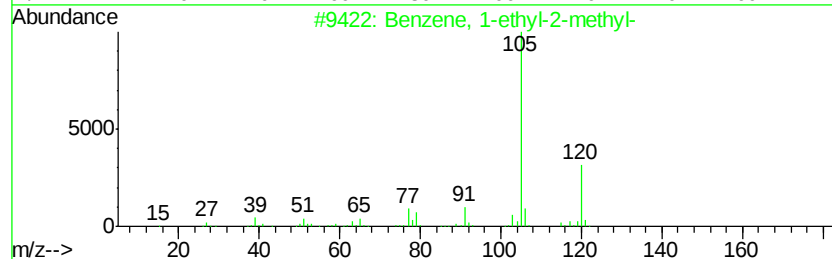
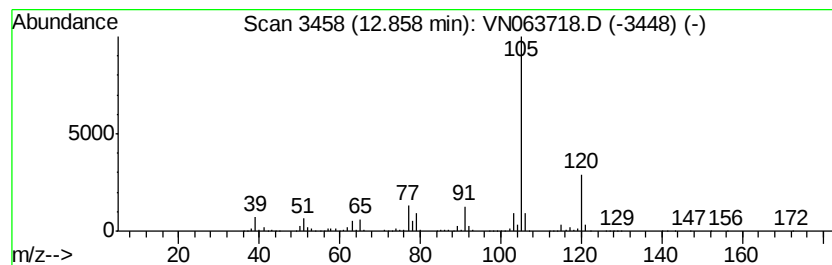
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
Quant Title : SW846 8260

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

Peak Number 1 Benzene, 1-ethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.86	167.49 ug/l	5067190	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
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, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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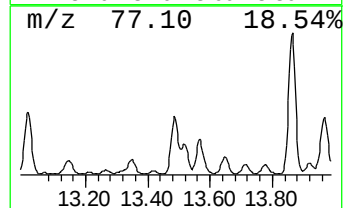
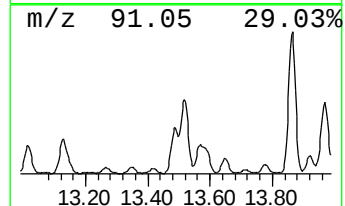
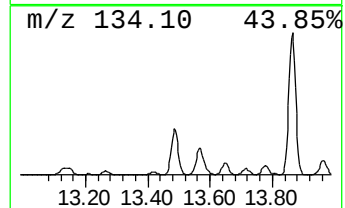
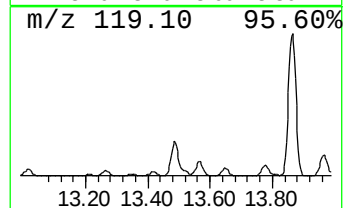
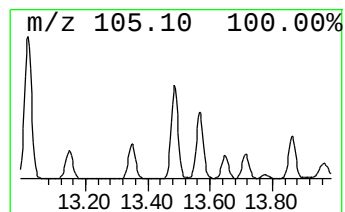
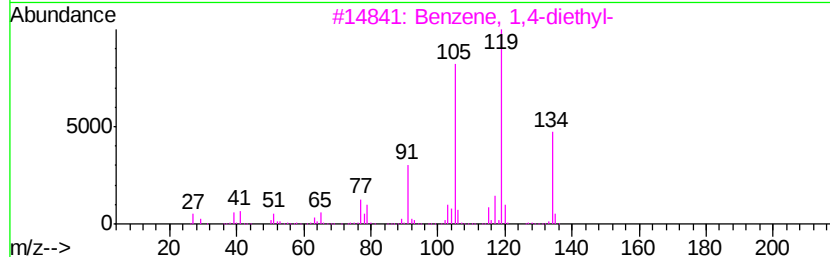
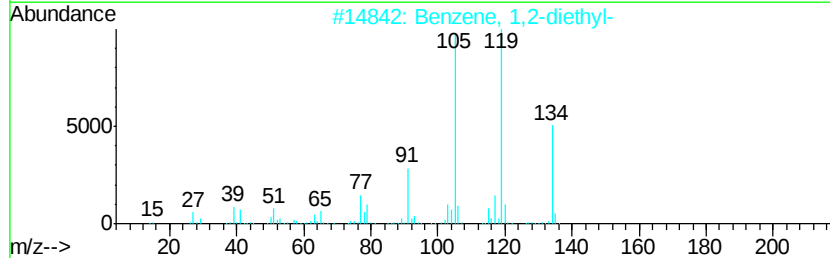
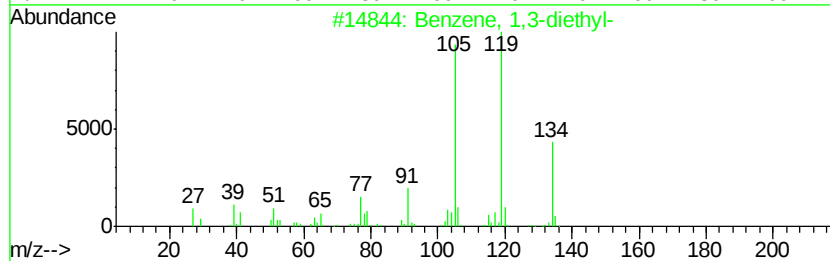
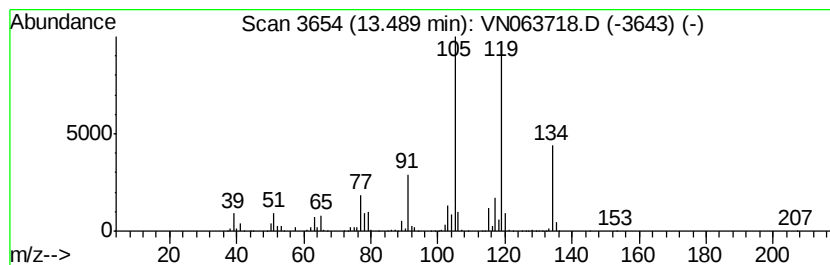
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Quant Title : SW846 8260

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

Peak Number 2 Benzene, 1,3-diethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.49	167.11 ug/l	5055760	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	97
2		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	96
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
5		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91



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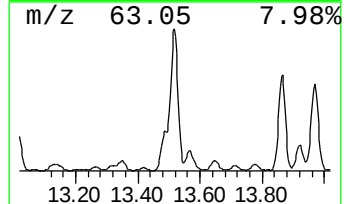
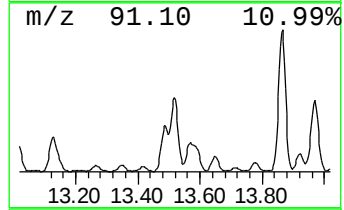
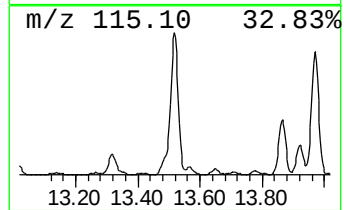
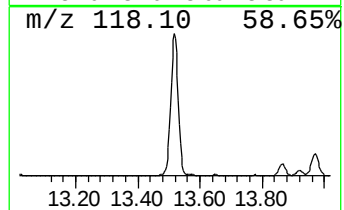
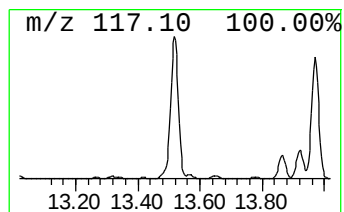
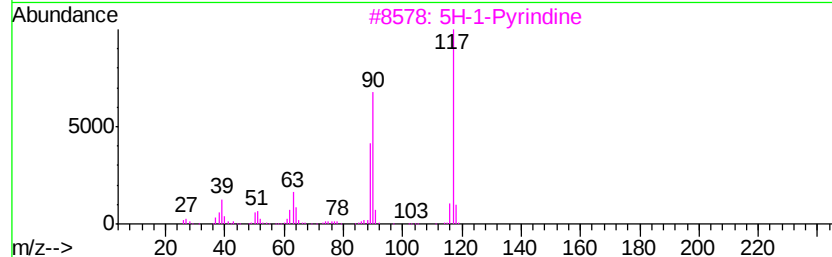
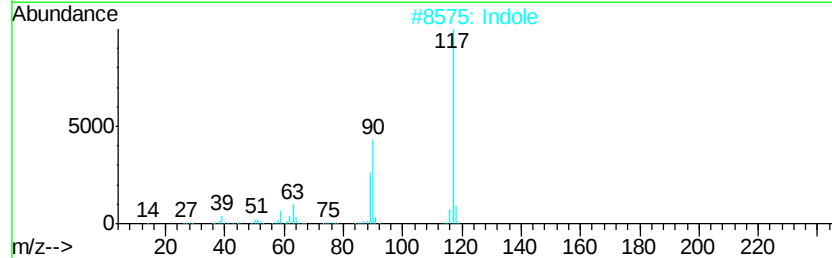
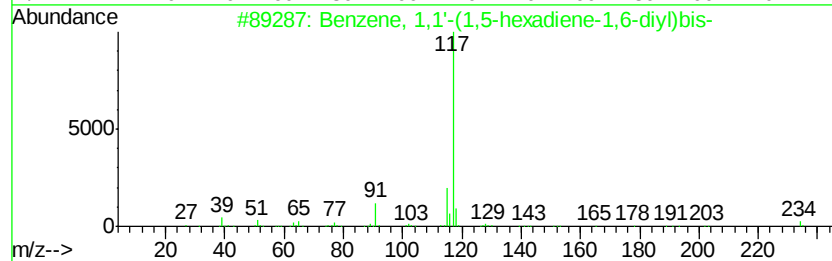
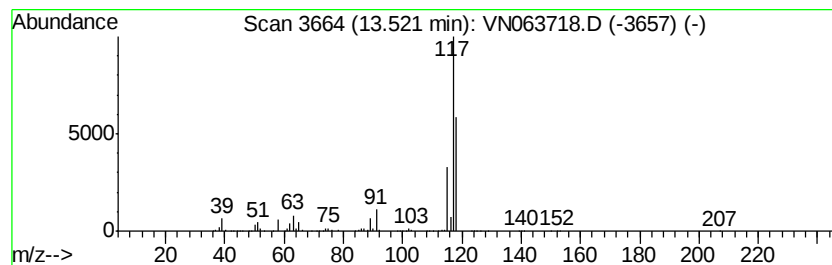
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Quant Title : SW846 8260

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

Peak Number 3 Benzene, 1,1'-(1,5-hexadien... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.52	328.80 ug/l	9947410	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	59
2		Indole	117	C8H7N	000120-72-9	43
3		5H-1-Pyrindine	117	C8H7N	000270-91-7	43
4		Benzaldehyde, 4-(1-phenyl-2-prop...	238	C16H14O2	1000277-56-1	42
5		Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	38



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092820\
Data File : VN063718.D
Acq On : 28 Sep 2020 20:20
Operator : JC/MD
Sample : L4178-06
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TP-6

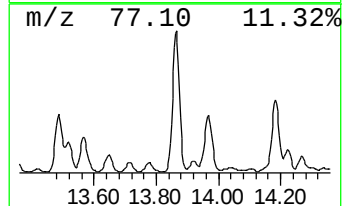
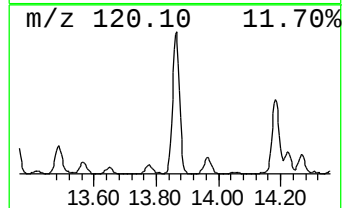
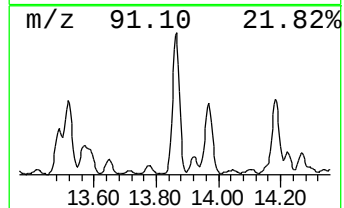
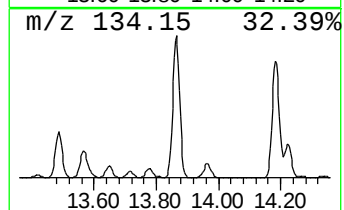
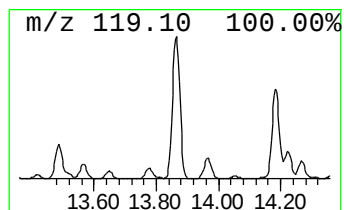
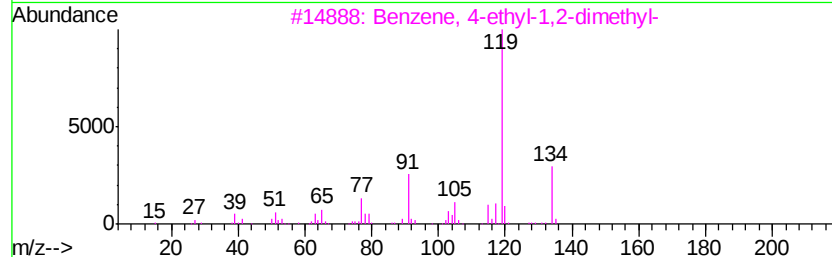
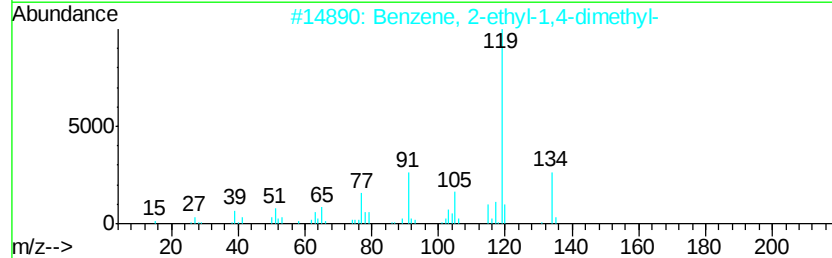
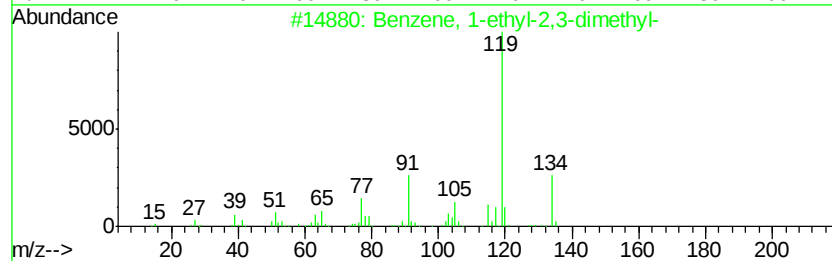
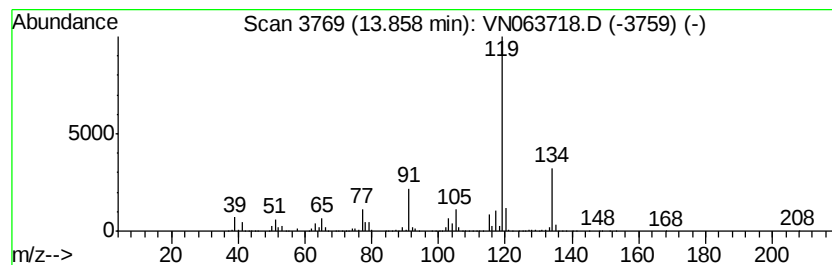
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Quant Title : SW846 8260

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

Peak Number 4 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.86	465.15 ug/l	14072500	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
2		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
4		o-Cymene	134	C10H14	000527-84-4	95
5		Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092820\
Data File : VN063718.D
Acq On : 28 Sep 2020 20:20
Operator : JC/MD
Sample : L4178-06
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampled :
TP-6

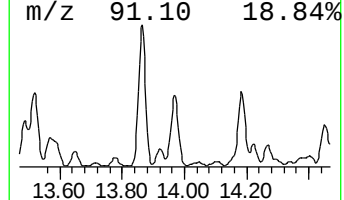
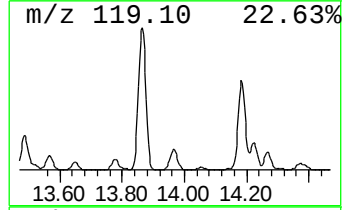
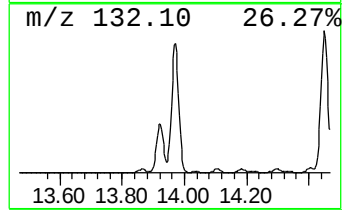
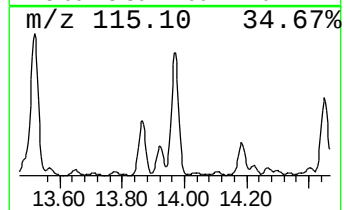
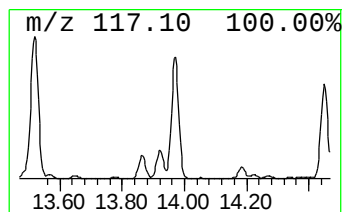
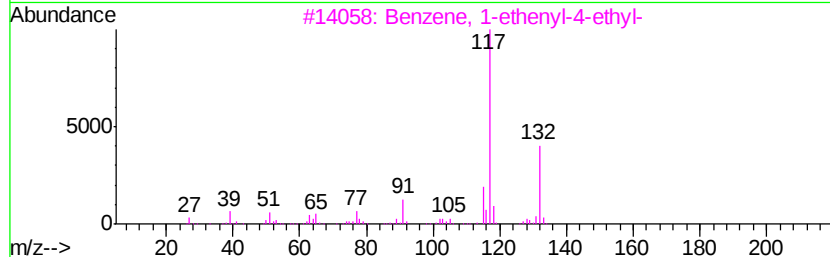
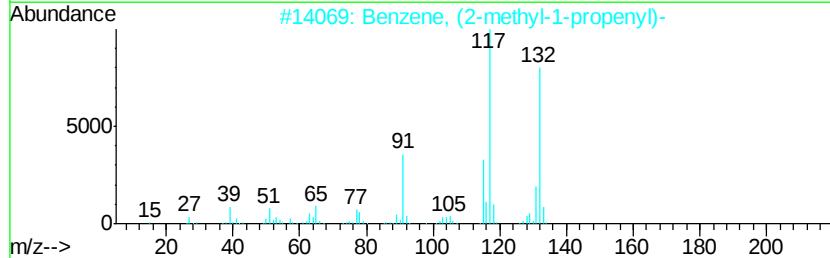
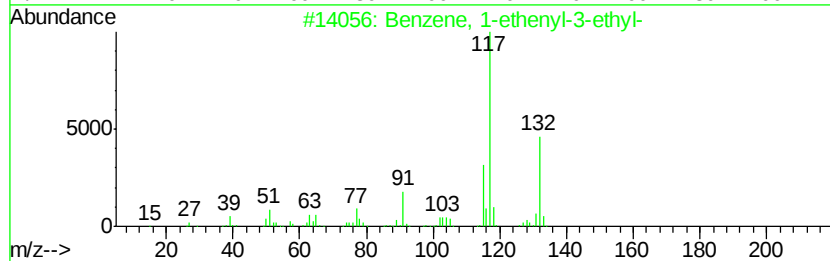
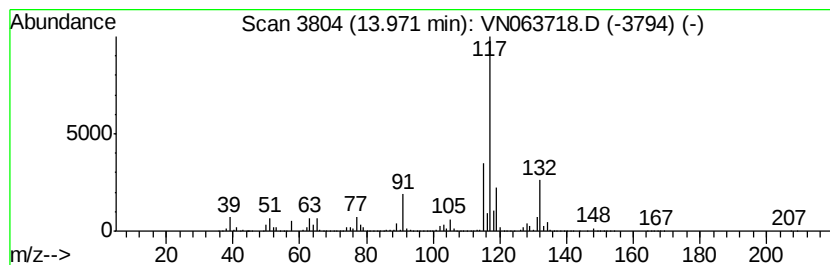
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Quant Title : SW846 8260

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

Peak Number 5 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.97	317.04 ug/l	9591650	1,4-Dichlorobenzene-d4	13.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
2		Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	81
3		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76
4		Indan, 1-methyl-	132	C10H12	000767-58-8	76
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092820\
Data File : VN063718.D
Acq On : 28 Sep 2020 20:20
Operator : JC/MD
Sample : L4178-06
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampled :
TP-6

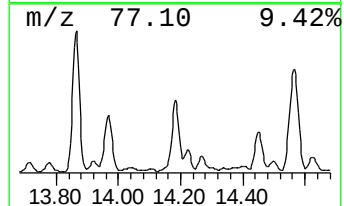
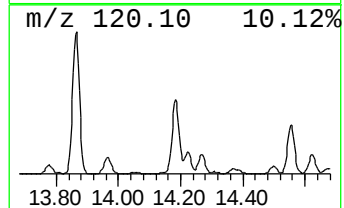
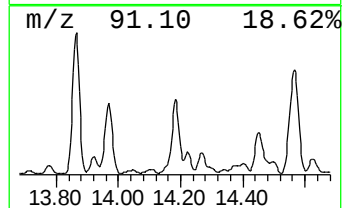
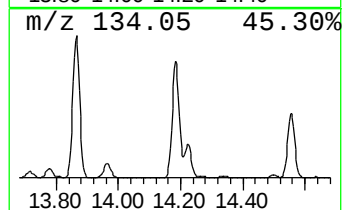
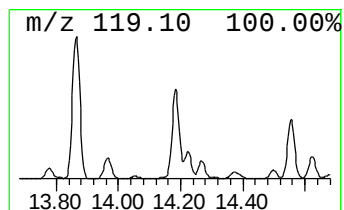
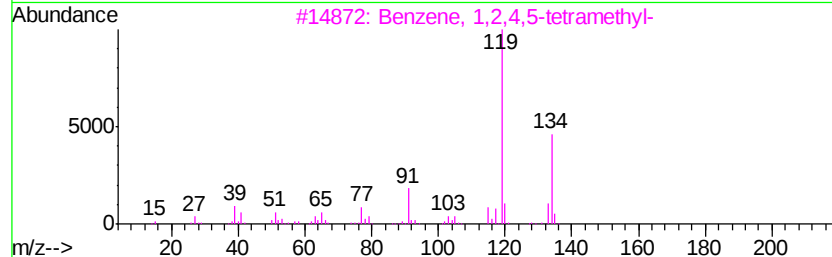
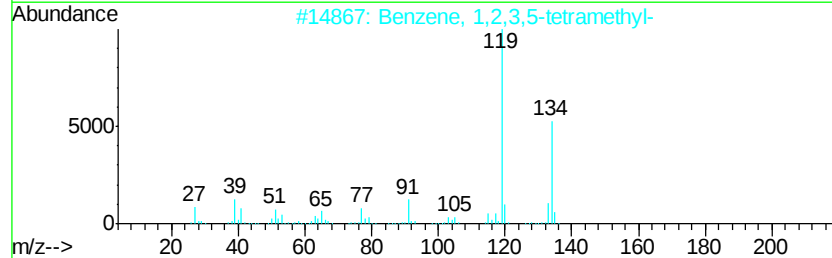
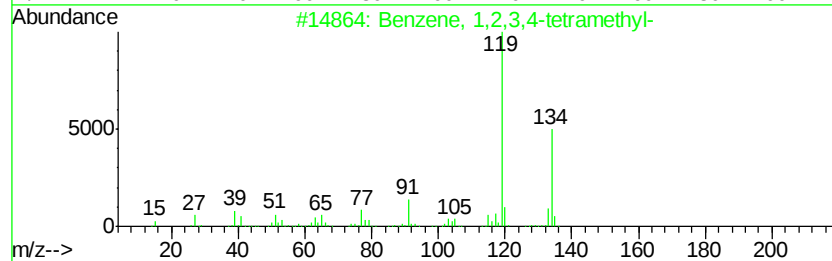
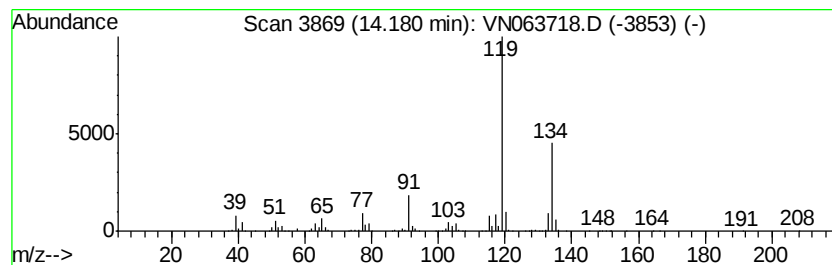
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.18	296.19 ug/l	8961020	1,4-Dichlorobenzene-d4	13.32

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	93



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092820\
Data File : VN063718.D
Acq On : 28 Sep 2020 20:20
Operator : JC/MD
Sample : L4178-06
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TP-6

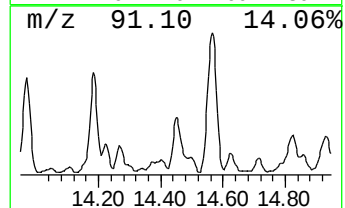
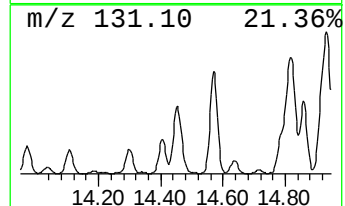
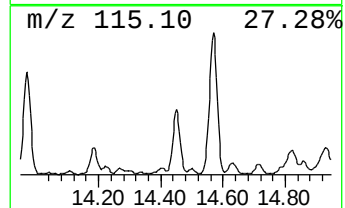
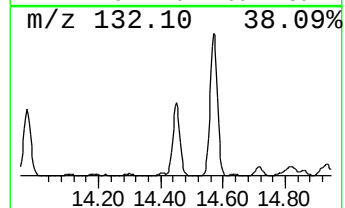
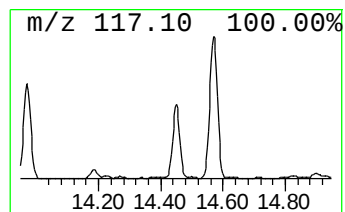
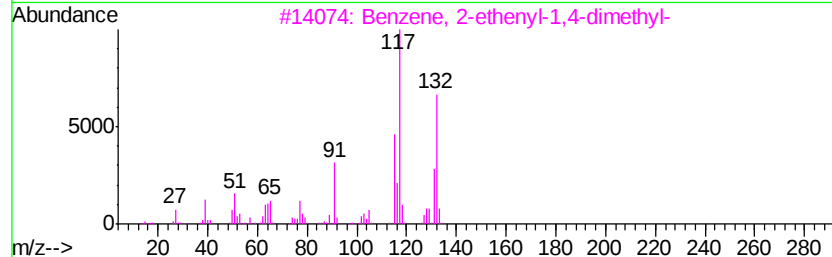
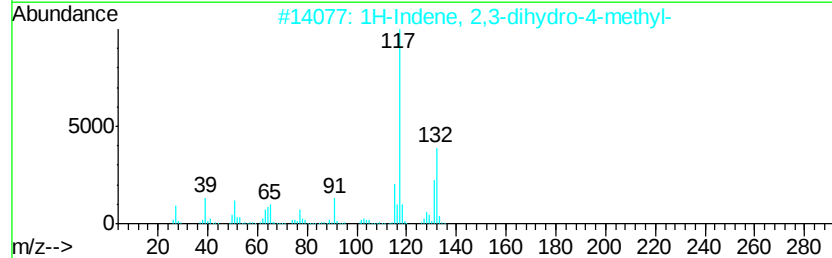
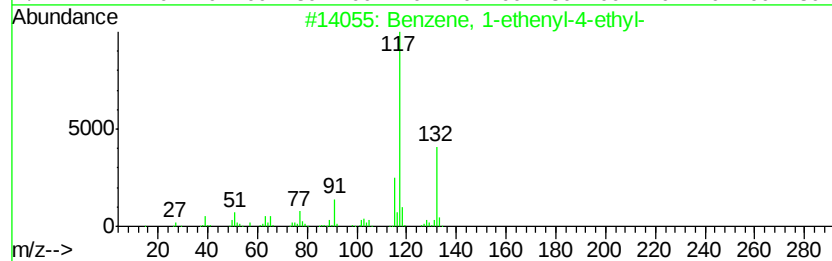
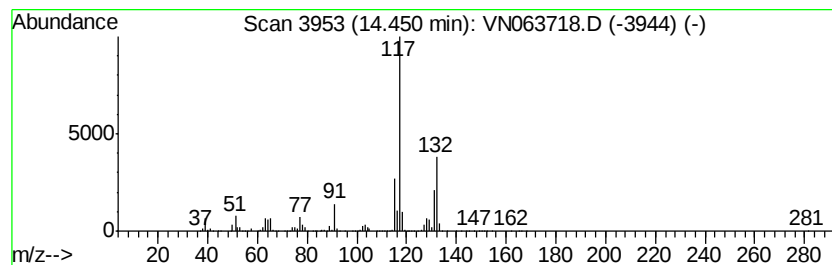
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Quant Title : SW846 8260

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

Peak Number 7 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.45	214.80 ug/l	6498670	1,4-Dichlorobenzene-d4	13.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
3		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	93
4		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	92
5		3-Phenylbut-1-ene	132	C10H12	000934-10-1	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092820\
Data File : VN063718.D
Acq On : 28 Sep 2020 20:20
Operator : JC/MD
Sample : L4178-06
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TP-6

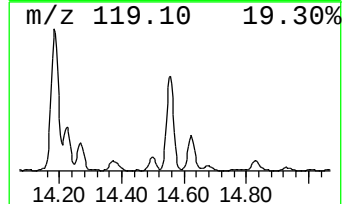
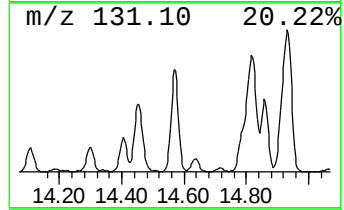
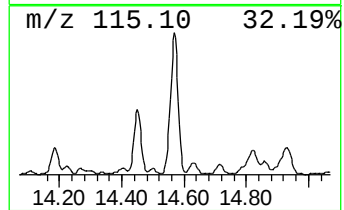
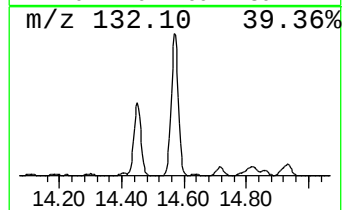
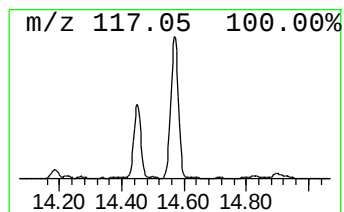
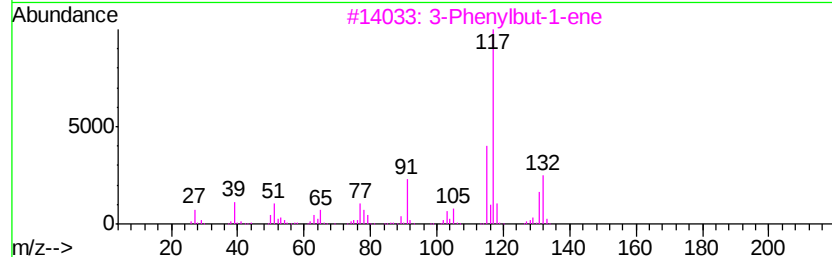
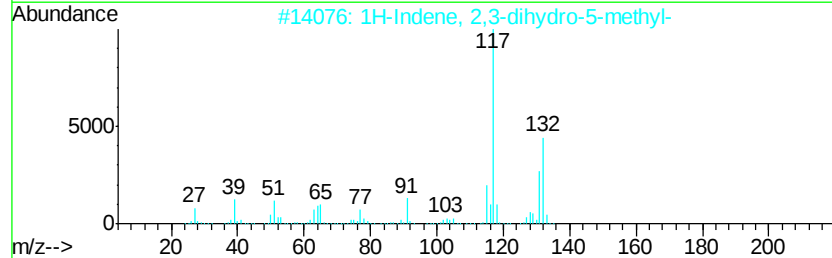
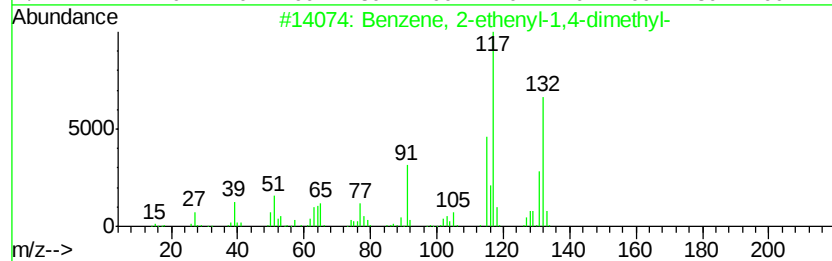
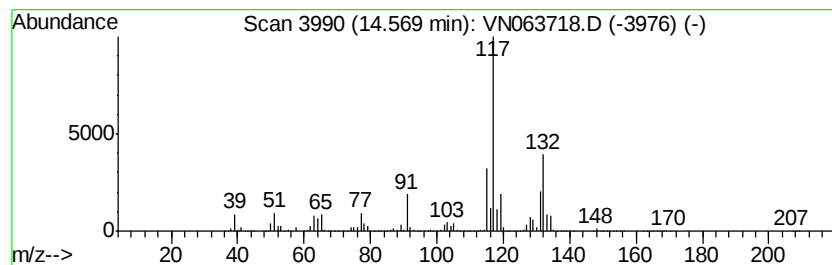
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Quant Title : SW846 8260

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

Peak Number 8 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	639.40 ug/l	19344300	1,4-Dichlorobenzene-d4	13.32

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	93
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, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	76



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092820\
Data File : VN063718.D
Acq On : 28 Sep 2020 20:20
Operator : JC/MD
Sample : L4178-06
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampled :
TP-6

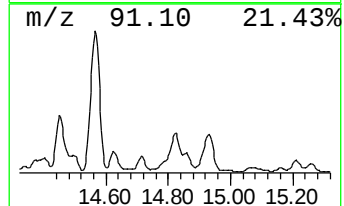
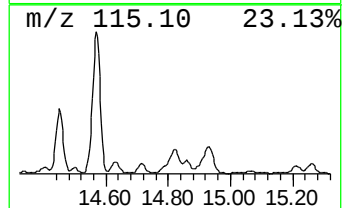
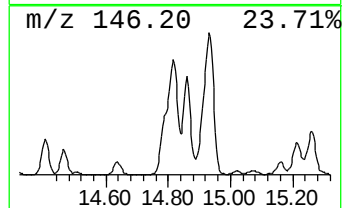
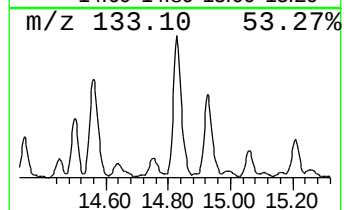
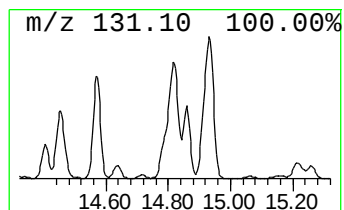
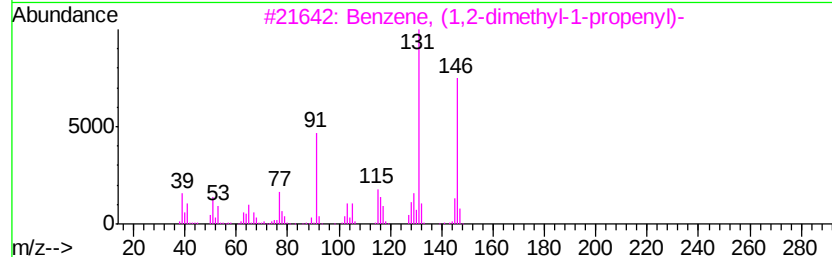
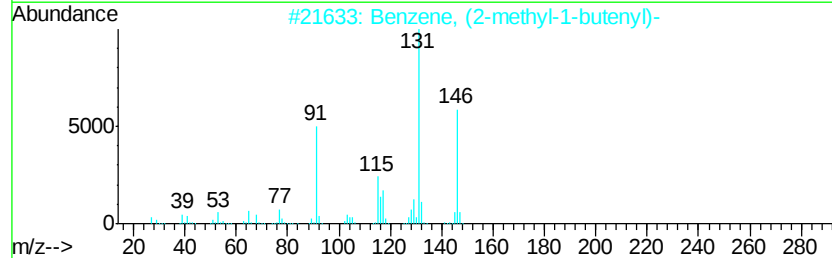
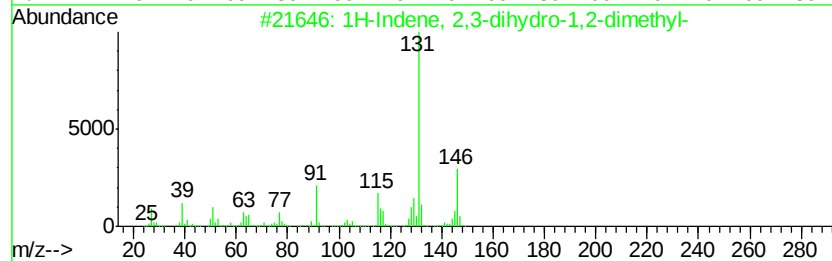
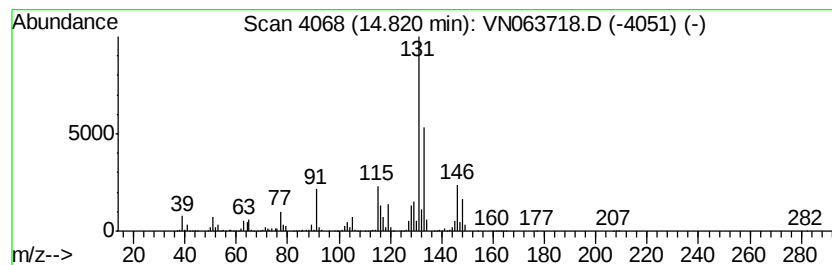
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Quant Title : SW846 8260

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

Peak Number 9 1H-Indene, 2,3-dihydro-1,2-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	180.46 ug/l	5459740	1,4-Dichlorobenzene-d4	13.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	92
2		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	91
3		Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	90
4		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	78
5		Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA N\DATA\VN092820\
Data File : VN063718.D
Acq On : 28 Sep 2020 20:20
Operator : JC/MD
Sample : L4178-06
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 22 Sample Multiplier: 1

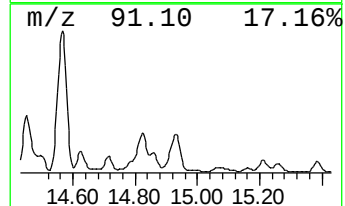
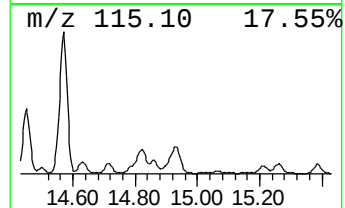
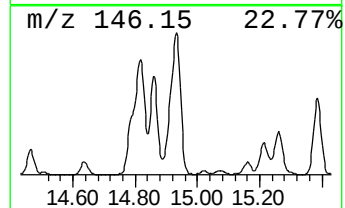
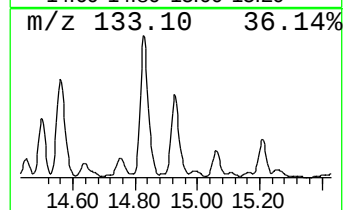
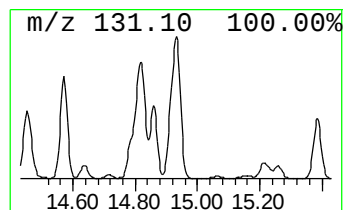
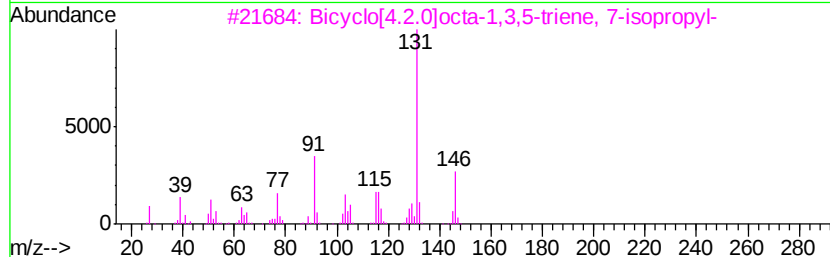
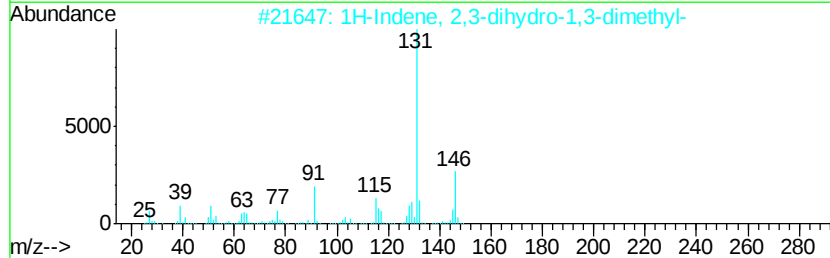
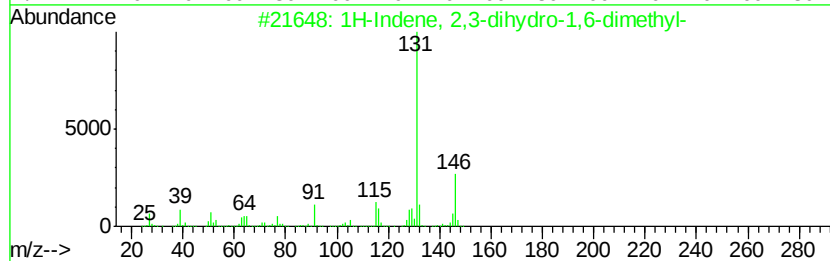
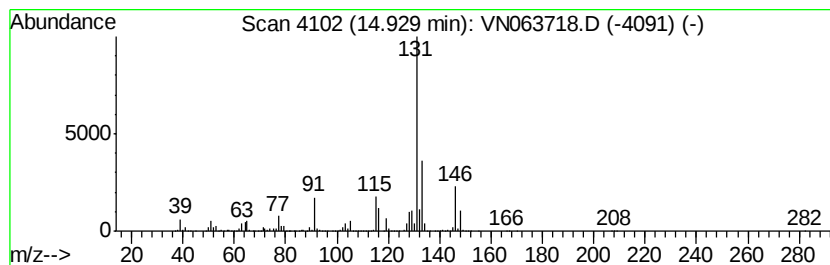
Instrument :
MSVOA_N
ClientSampled :
TP-6

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.M
Quant Title : SW846 8260

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

Peak Number 10 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.93	181.40 ug/l	5488130	1,4-Dichlorobenzene-d4	13.32	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	81
2	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	81
3	Bicyclo[4.2.0]octa-1,3,5-triene...	146	C11H14	027087-54-3	81
4	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	81
5	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	81



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN092820\
Data File : VN063718.D
Acq On : 28 Sep 2020 20:20
Operator : JC/MD
Sample : L4178-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 22 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
TP-6

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N091620W.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-ethyl-...	12.86	167.5	ug/l	5067190	4	13.32	1512700	50.0
Benzene, 1,3-diet...	13.49	167.1	ug/l	5055760	4	13.32	1512700	50.0
Benzene, 1,1'-(1,...	13.52	328.8	ug/l	9947410	4	13.32	1512700	50.0
Benzene, 1-ethyl-...	13.86	465.1	ug/l	14072500	4	13.32	1512700	50.0
Benzene, 1-etheny...	13.97	317.0	ug/l	9591650	4	13.32	1512700	50.0
Benzene, 1,2,3,4-...	14.18	296.2	ug/l	8961020	4	13.32	1512700	50.0
Benzene, 1-etheny...	14.45	214.8	ug/l	6498670	4	13.32	1512700	50.0
Benzene, 2-etheny...	14.57	639.4	ug/l	19344300	4	13.32	1512700	50.0
1H-Indene, 2,3-di...	14.82	180.5	ug/l	5459740	4	13.32	1512700	50.0
1H-Indene, 2,3-di...	14.93	181.4	ug/l	5488130	4	13.32	1512700	50.0