

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_Y\DATA\VY110719\
 Data File : VY000603.D
 Acq On : 08 Nov 2019 00:37
 Operator : SY/MD
 Sample : K5722-07
 Misc : 7.36G/5ML/MSVOA_Y/SOIL
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 5-C

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_Y\METHODS\82Y103019S.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	3.127	223	228	237	rVB3	9112	20922	1.63%	0.260%
2	4.725	480	490	504	rBV3	34374	100467	7.84%	1.248%
3	5.755	647	659	674	rBV4	30425	98277	7.67%	1.220%
4	7.730	973	983	989	rBV2	157071	379378	29.59%	4.711%
5	7.803	989	995	1007	rVB2	301231	672108	52.43%	8.347%
6	8.157	1041	1053	1065	rBV	174304	384739	30.01%	4.778%
7	8.699	1134	1142	1152	rBV	487438	945527	73.76%	11.742%
8	10.181	1378	1385	1393	rBV	741897	1281982	100.00%	15.920%
9	11.492	1592	1600	1609	rVB	707479	1130680	88.20%	14.041%
10	12.479	1750	1762	1778	rBV	547864	916741	71.51%	11.385%
11	13.321	1895	1900	1901	rBV4	12541	16447	1.28%	0.204%
12	13.424	1911	1917	1925	rVB	692575	1060557	82.73%	13.171%
13	13.875	1985	1991	1995	rBV3	12420	21955	1.71%	0.273%
14	14.314	2058	2063	2071	rBV5	11292	22927	1.79%	0.285%
15	14.674	2115	2122	2131	rBV4	19247	58333	4.55%	0.724%
16	14.869	2146	2154	2159	rBV8	9812	27666	2.16%	0.344%
17	15.095	2186	2191	2203	rVB5	36198	109412	8.53%	1.359%
18	15.235	2203	2214	2229	rBV3	52199	165957	12.95%	2.061%
19	15.540	2253	2264	2277	rBV6	33426	128501	10.02%	1.596%
20	15.729	2283	2295	2303	rBV6	37848	124596	9.72%	1.547%
21	15.808	2303	2308	2317	rVB6	10039	25605	2.00%	0.318%
22	16.040	2337	2346	2359	rVB5	25871	112372	8.77%	1.396%
23	16.479	2414	2418	2423	rVB7	10658	22554	1.76%	0.280%
24	16.570	2423	2433	2436	rBV3	23215	61692	4.81%	0.766%
25	16.936	2482	2493	2501	rBV8	38741	163043	12.72%	2.025%

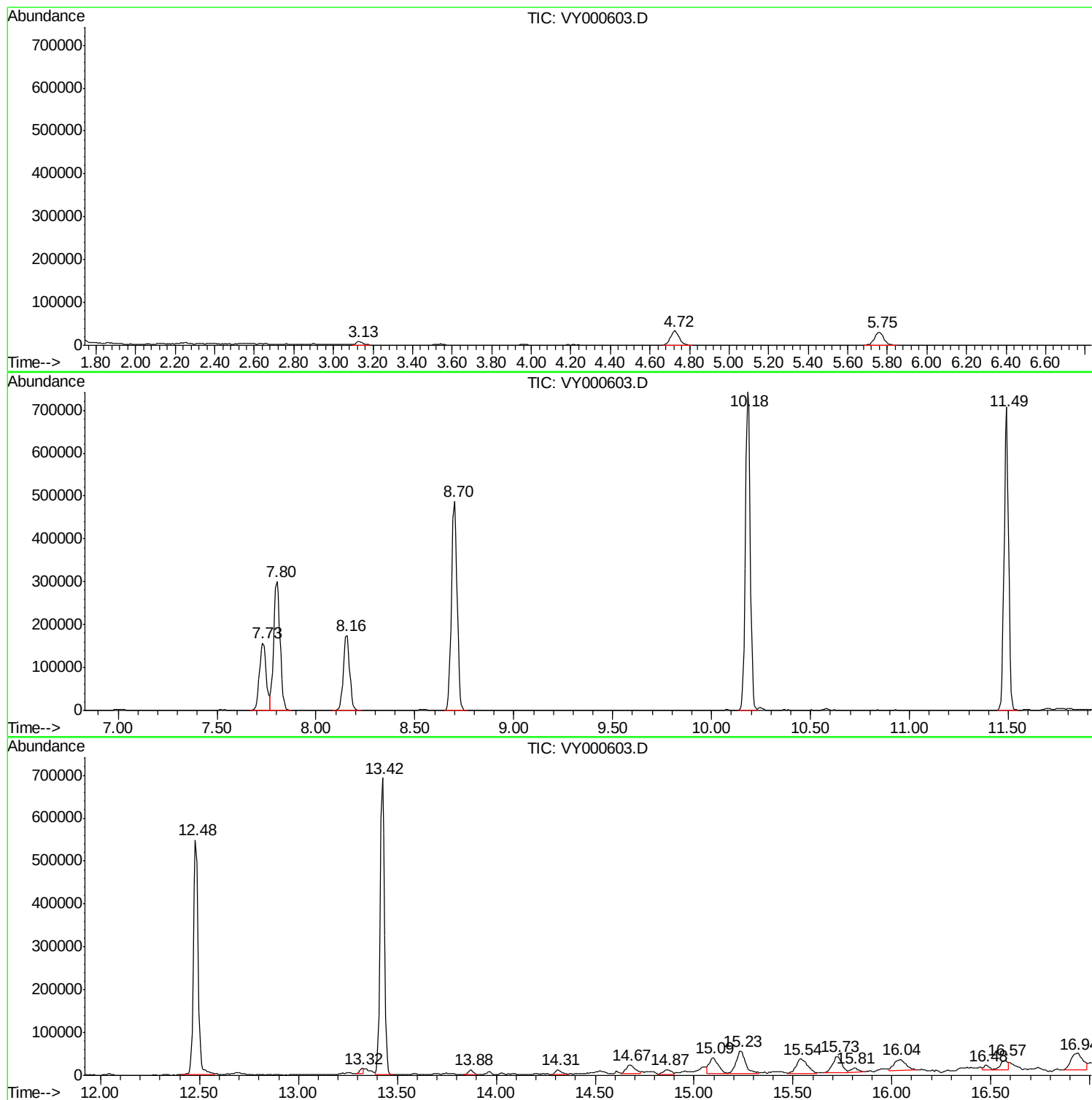
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Quant Method : Z:\V0ASRV\HPCHEM1\MSV0A_Y\METHODS\82Y103019S.M
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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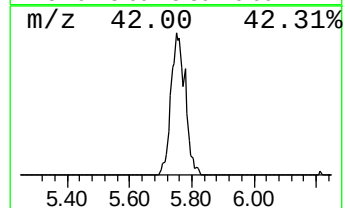
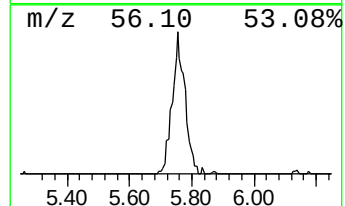
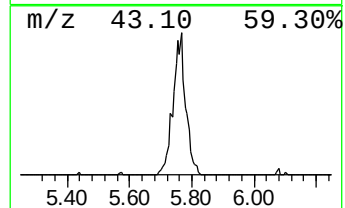
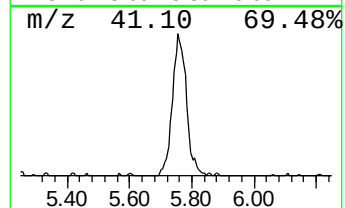
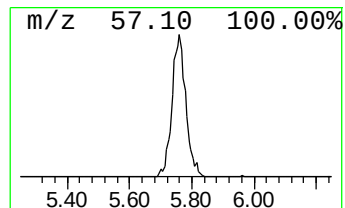
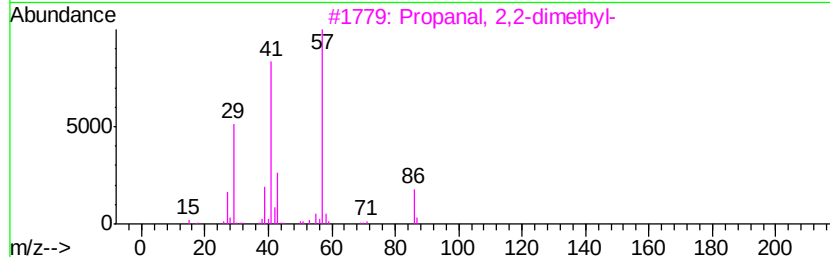
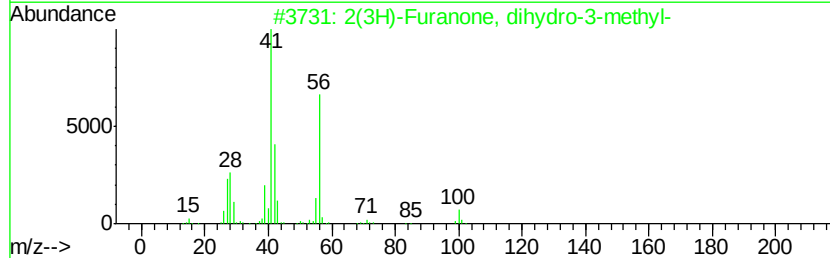
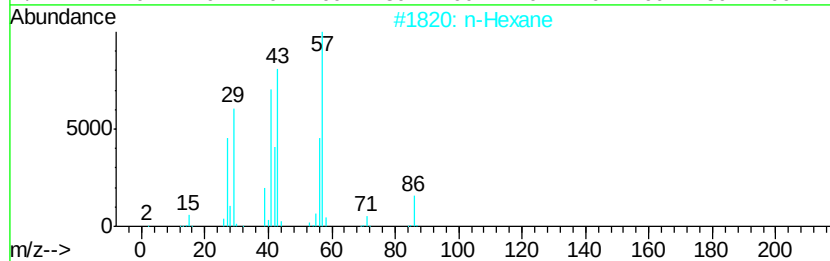
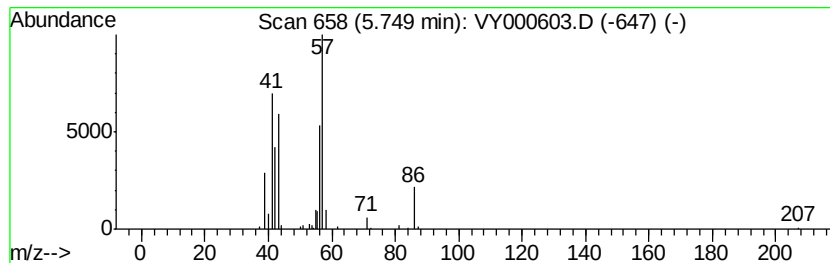
Quant Method : Z:\V0ASRV\HPCHEM1\MSV0A_Y\METHODS\82Y103019S.M
Quant Title : SW846 8260

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

Peak Number 1 n-Hexane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.75	7.31 ug/l	98277	Pentafluorobenzene	7.80

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexane	86	C6H14	000110-54-3	86
2			2(3H)-Furanone, dihydro-3-methyl-	100	C5H8O2	001679-47-6	53
3			Propanal, 2,2-dimethyl-	86	C5H10O	000630-19-3	43
4			1-Pentene, 4-methyl-	84	C6H12	000691-37-2	42
5			Pentane, 3-methyl-	86	C6H14	000096-14-0	40



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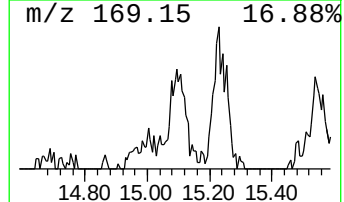
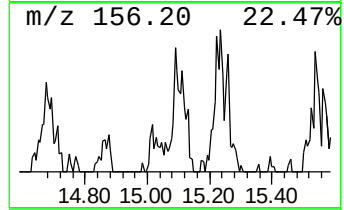
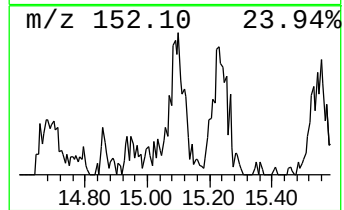
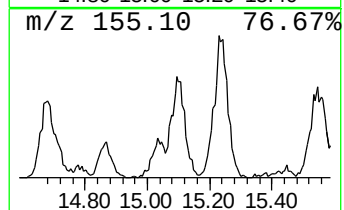
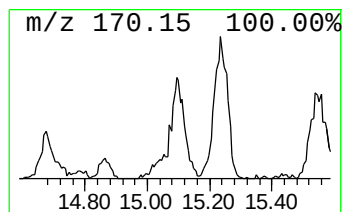
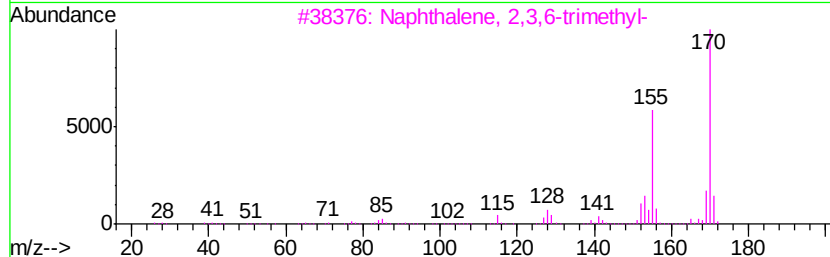
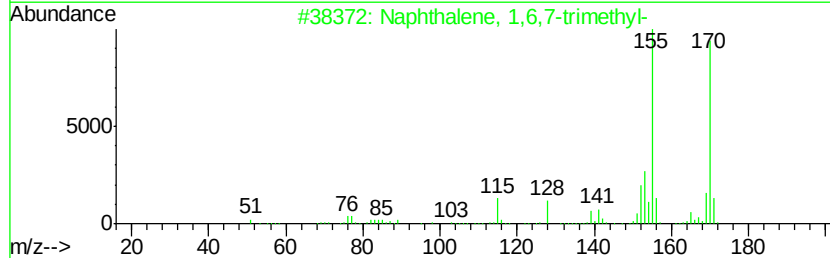
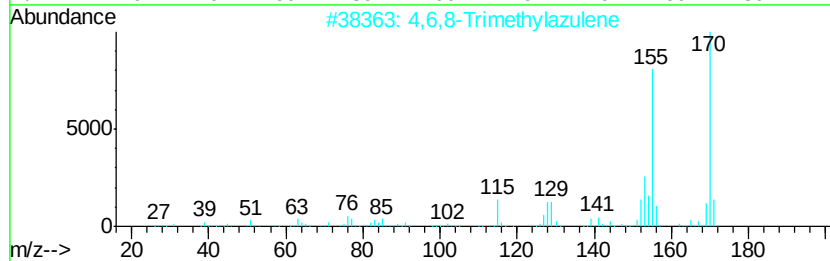
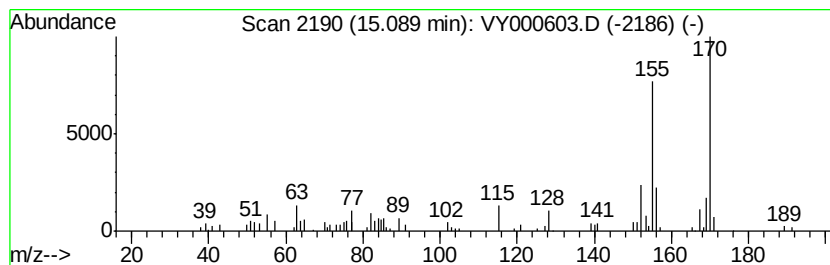
Instrument :
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ClientSampleId :
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Quant Title : SW846 8260

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Peak Number 2 4,6,8-Trimethylazulene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.09	5.16 ug/l	109412	1,4-Dichlorobenzene-d4	13.42
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	4,6,8-Trimethylazulene	170 C13H14	000941-81-1	95
2	Naphthalene, 1,6,7-trimethyl-	170 C13H14	002245-38-7	95
3	Naphthalene, 2,3,6-trimethyl-	170 C13H14	000829-26-5	94
4	Naphthalene, 1,4,5-trimethyl-	170 C13H14	002131-41-1	93
5	Naphthalene, 1,4,6-trimethyl-	170 C13H14	002131-42-2	91



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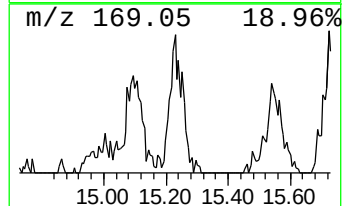
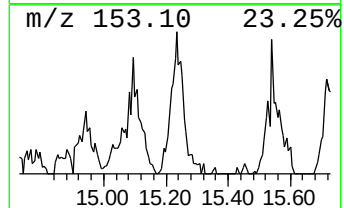
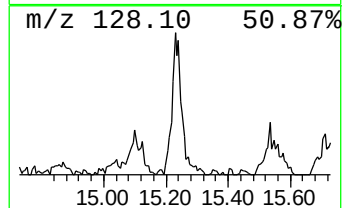
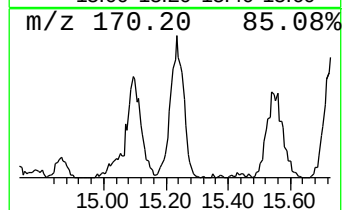
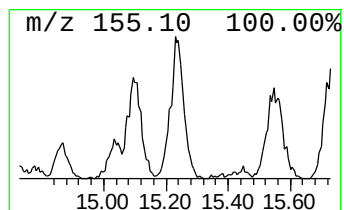
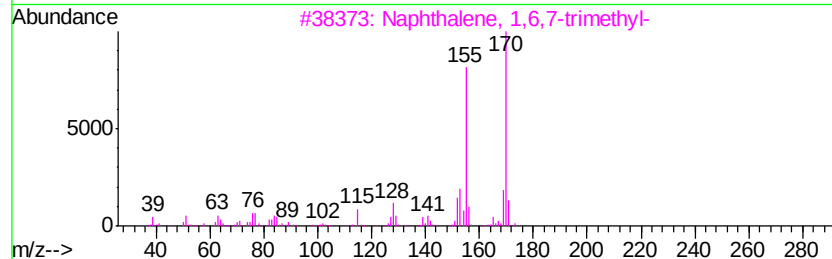
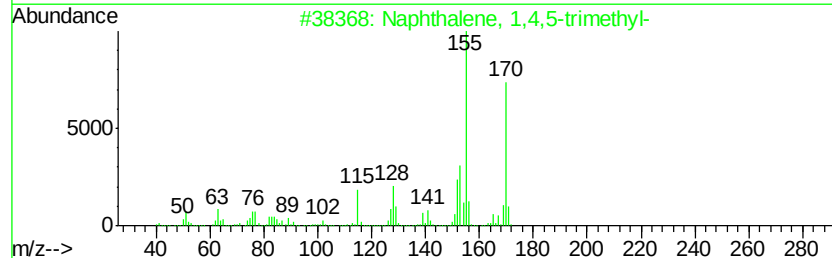
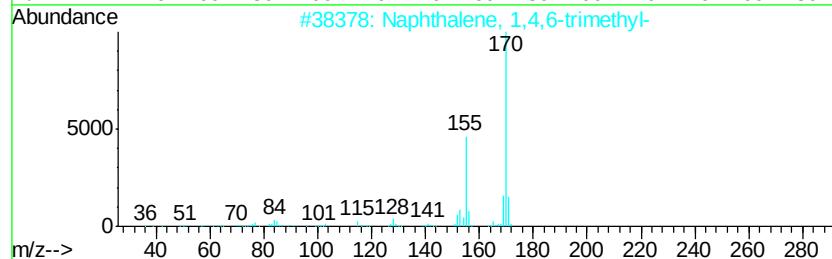
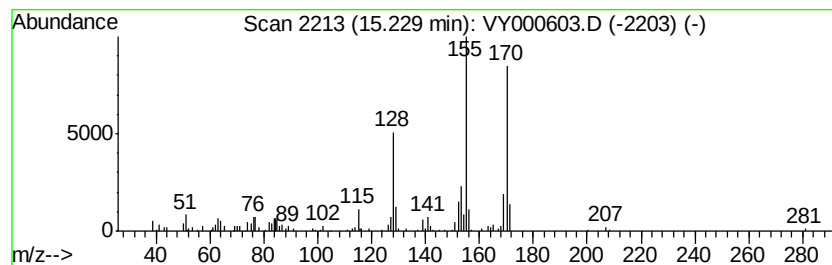
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TIC Integration Parameters: LSCINT.P

Peak Number 3 Naphthalene, 1,4,6-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.23	7.82 ug/l	165957	1,4-Dichlorobenzene-d4	13.42

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	93
2		Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	90
3		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	89
4		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	76
5		4,6,8-Trimethylazulene	170	C13H14	000941-81-1	72



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ALS Vial : 41 Sample Multiplier: 1

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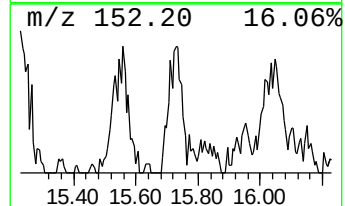
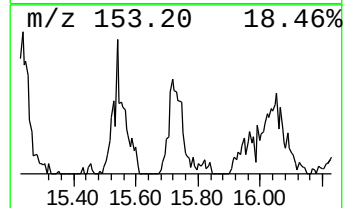
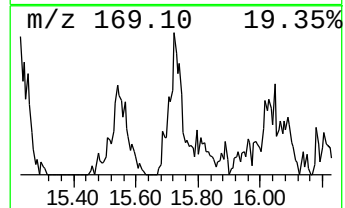
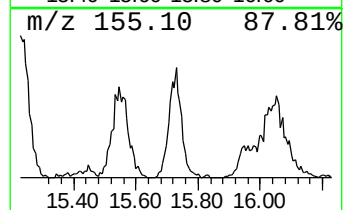
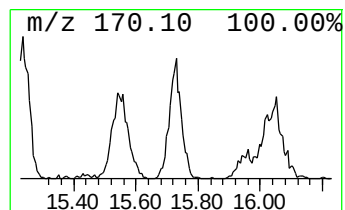
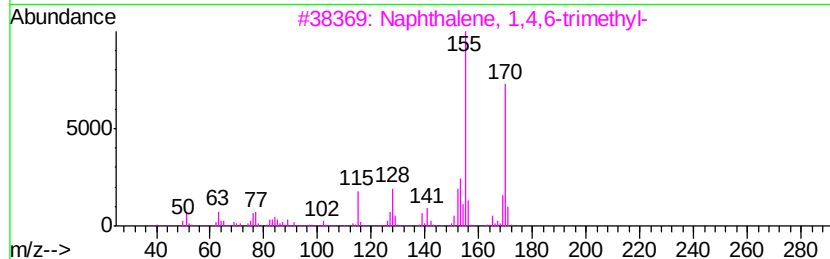
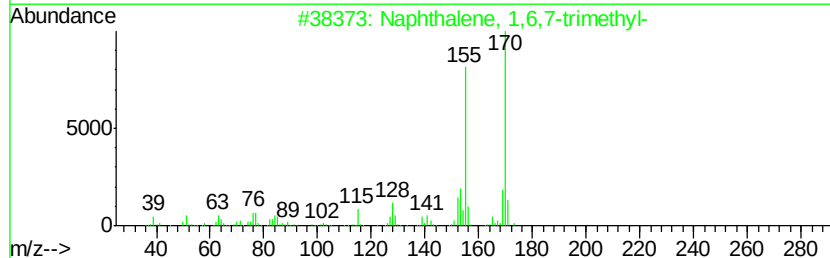
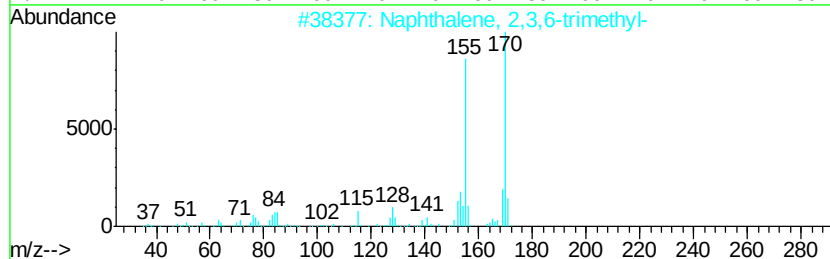
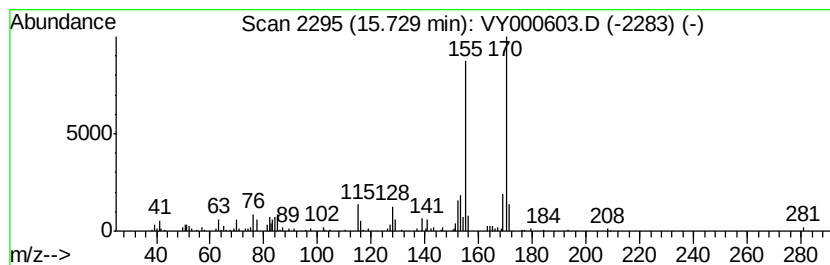
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Naphthalene, 2,3,6-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.73	5.87 ug/l	124596	1,4-Dichlorobenzene-d4	13.42

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
2		Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
3		Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	95
4		4,6,8-Trimethylazulene	170	C13H14	000941-81-1	80
5		Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	80



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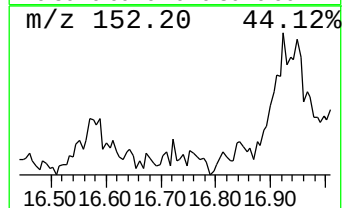
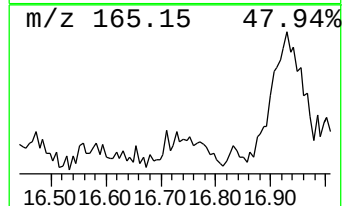
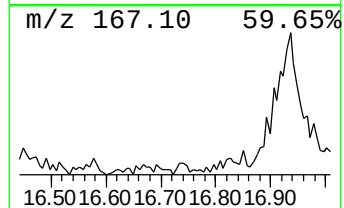
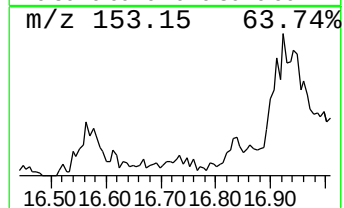
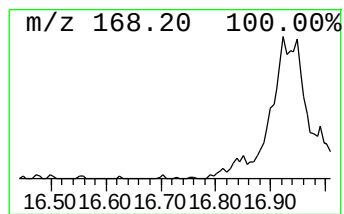
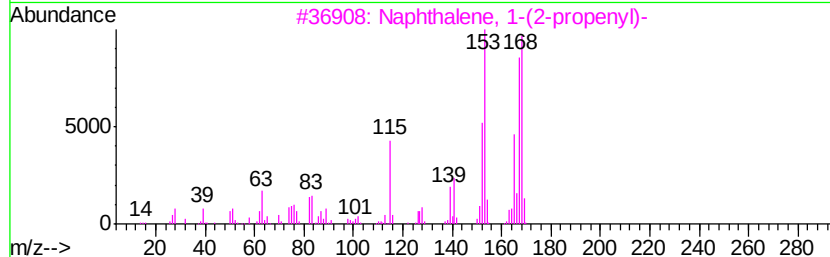
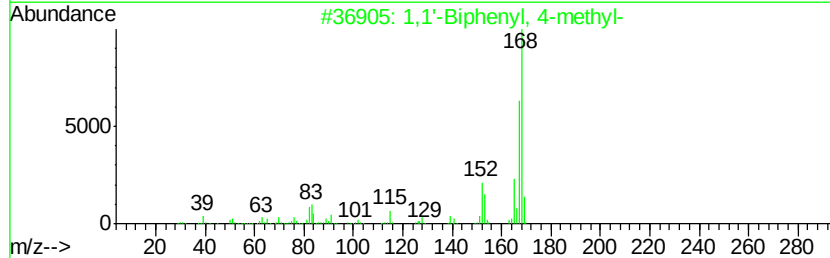
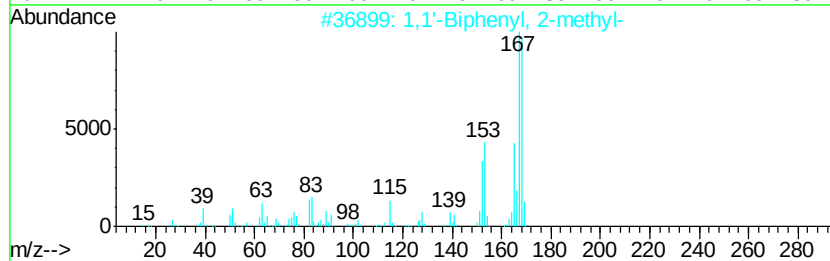
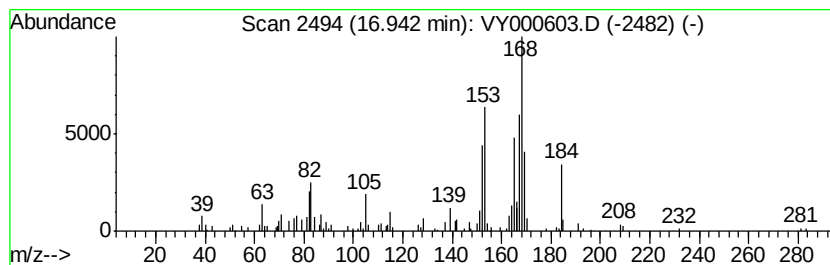
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 1,1'-Biphenyl, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.94	7.69 ug/l	163043	1,4-Dichlorobenzene-d4	13.42	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2-methyl-	168	C13H12	000643-58-3	83
2	1,1'-Biphenyl, 4-methyl-	168	C13H12	000644-08-6	70
3	Naphthalene, 1-(2-propenyl)-	168	C13H12	002489-86-3	53
4	1-Isopropenylnaphthalene	168	C13H12	001855-47-6	49
5	Diphenylmethane	168	C13H12	000101-81-5	47



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
n-Hexane	5.75	7.3	ug/l	98277	1	7.80	672108	50.0
4,6,8-Trimethylaz...	15.09	5.2	ug/l	109412	4	13.42	1060560	50.0
Naphthalene, 1,4,...	15.23	7.8	ug/l	165957	4	13.42	1060560	50.0
Naphthalene, 2,3,...	15.73	5.9	ug/l	124596	4	13.42	1060560	50.0
1,1'-Biphenyl, 2-...	16.94	7.7	ug/l	163043	4	13.42	1060560	50.0