

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX091818\
 Data File : VX004650.D
 Acq On : 18 Sep 2018 11:49
 Operator : JC/MD
 Sample : J4898-01 50X
 Misc : 5.26g/5mL/100uL/5.0mL/MSVOA_X/MEOH
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 ROLLOFF01-COMP-FILL

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_X\METHOD\82X091118W.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	5.501	701	713	728	rBV2	144026	408198	6.97%	2.323%
2	5.659	728	739	755	rVV	196934	550497	9.40%	3.133%
3	6.068	793	806	813	rBV	139351	361544	6.18%	2.057%
4	6.141	813	818	832	rVB2	65019	170279	2.91%	0.969%
5	6.860	925	936	953	rBV	305392	716451	12.24%	4.077%
6	8.714	1233	1240	1247	rBV	672164	1061742	18.14%	6.042%
7	8.787	1247	1252	1265	rVB	236703	369549	6.31%	2.103%
8	10.116	1464	1470	1479	rBV	598945	807131	13.79%	4.593%
9	10.256	1487	1493	1502	rVB	126367	170560	2.91%	0.971%
10	10.360	1502	1510	1516	rBV	238059	321105	5.49%	1.827%
11	10.701	1561	1566	1576	rVB	90735	116432	1.99%	0.663%
12	11.140	1632	1638	1649	rBV	620085	767846	13.12%	4.369%
13	11.506	1693	1698	1710	rVB	50112	75215	1.29%	0.428%
14	11.811	1743	1748	1752	rVB	108780	128489	2.20%	0.731%
15	12.018	1776	1782	1787	rBV	86512	113734	1.94%	0.647%
16	12.079	1787	1792	1799	rVV	810146	985002	16.83%	5.605%
17	12.469	1847	1856	1862	rBV	614195	756915	12.93%	4.307%
18	13.024	1942	1947	1951	rBV	93920	134851	2.30%	0.767%
19	13.402	2004	2009	2014	rBV	95420	110785	1.89%	0.630%
20	13.481	2017	2022	2027	rVB	100275	132981	2.27%	0.757%
21	13.835	2073	2080	2087	rBV	4814348	5853286	100.00%	33.308%
22	13.932	2091	2096	2102	rVB	115699	166900	2.85%	0.950%
23	14.700	2217	2222	2232	rVB	989547	1246799	21.30%	7.095%
24	14.840	2239	2245	2257	rVB	548573	691244	11.81%	3.933%
25	15.273	2306	2316	2321	rBV	111793	155067	2.65%	0.882%
26	15.554	2355	2362	2373	rBV	136064	235502	4.02%	1.340%
27	15.688	2373	2384	2388	rVV	158671	277412	4.74%	1.579%
28	15.731	2388	2391	2398	rVB	100108	145511	2.49%	0.828%
29	15.920	2412	2422	2433	rVB2	50440	129927	2.22%	0.739%
30	16.176	2456	2464	2472	rBV	156074	282940	4.83%	1.610%
31	16.419	2496	2504	2512	rBV2	63476	129558	2.21%	0.737%

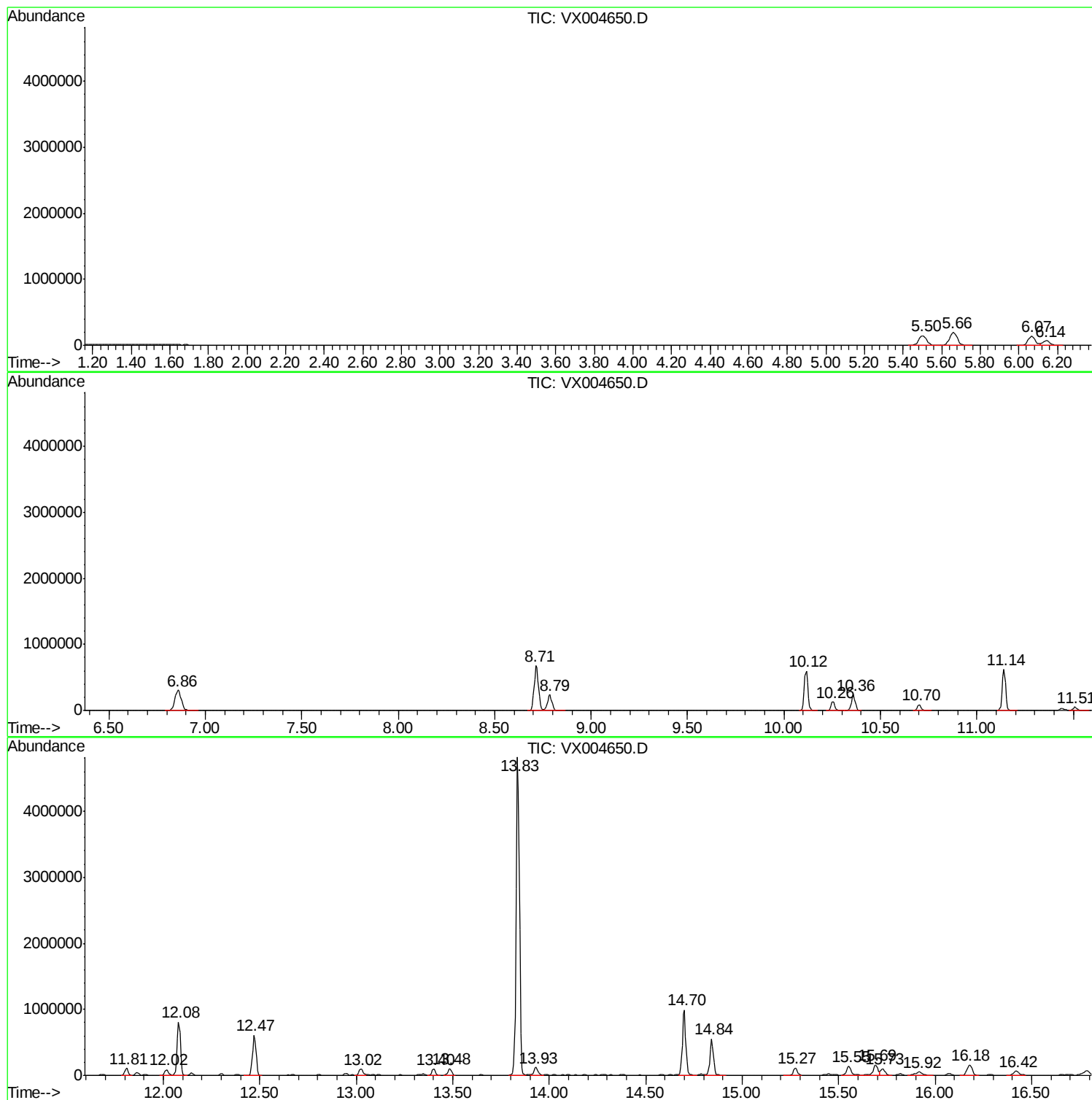
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ClientSampleId :
ROLLOFF01-COMP-FILL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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



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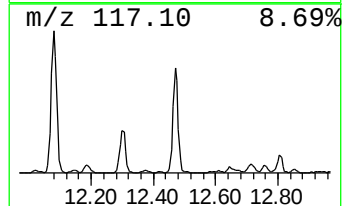
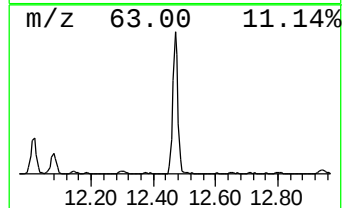
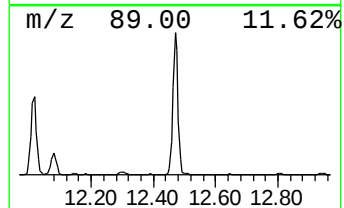
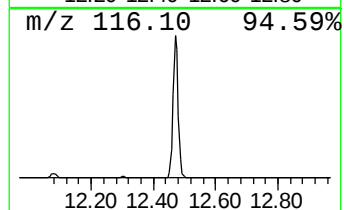
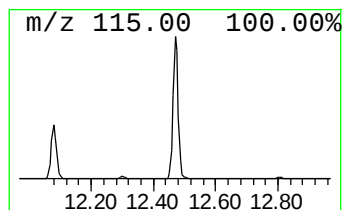
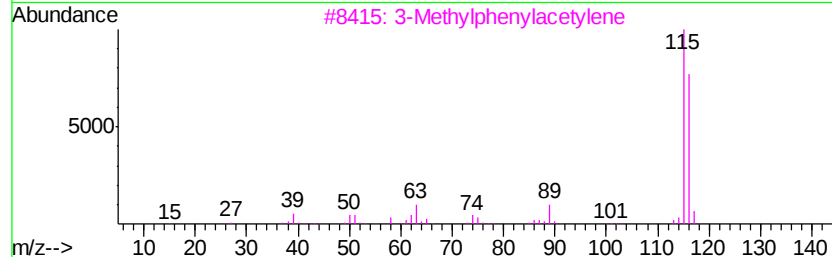
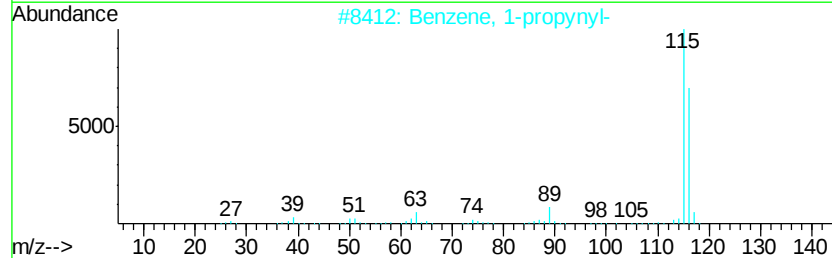
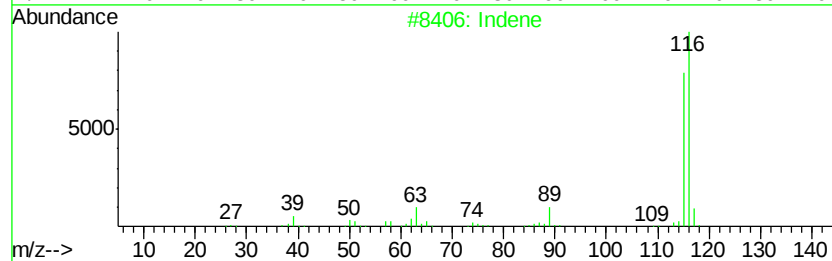
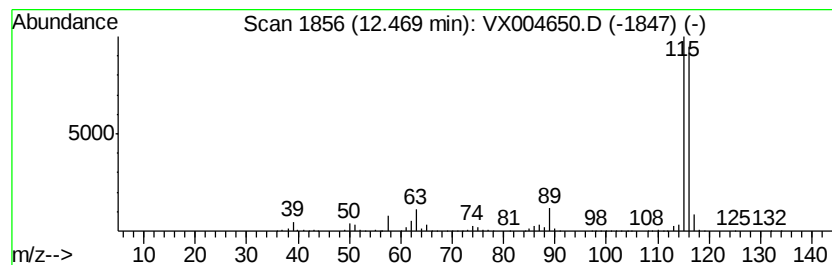
Instrument :
MSVOA_X
ClientSampleId :
ROLLOFF01-COMP-FILL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 1 Indene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.47	38.42 ug/l	756915	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Indene		116 C9H8	000095-13-6 97
2	Benzene, 1-propynyl-		116 C9H8	000673-32-5 94
3	3-Methylphenylacetylene		116 C9H8	000766-82-5 91
4	Benzene, 1-ethynyl-4-methyl-		116 C9H8	000766-97-2 91
5	2-Methylphenylacetylene		116 C9H8	000766-47-2 91



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ClientSampleId :
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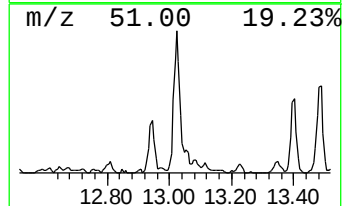
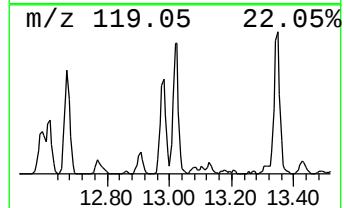
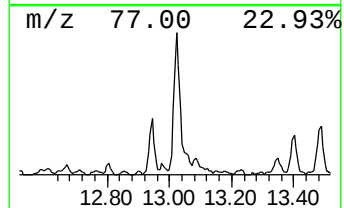
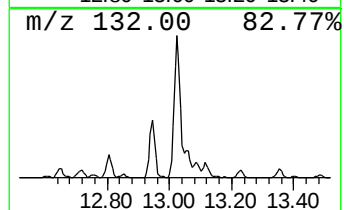
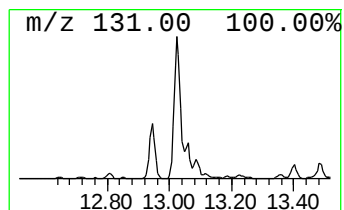
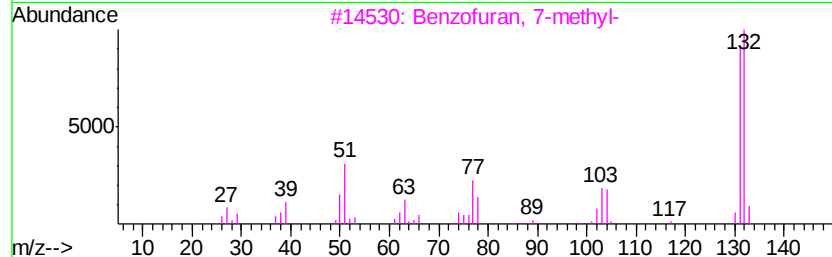
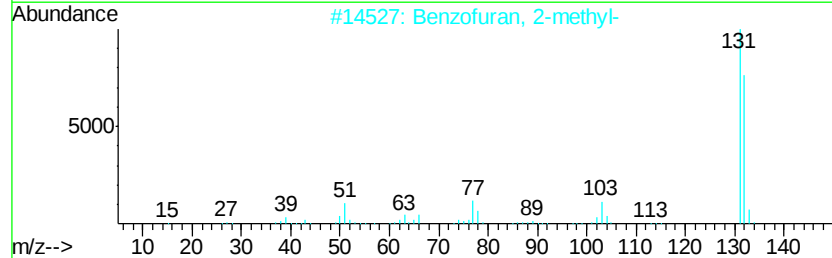
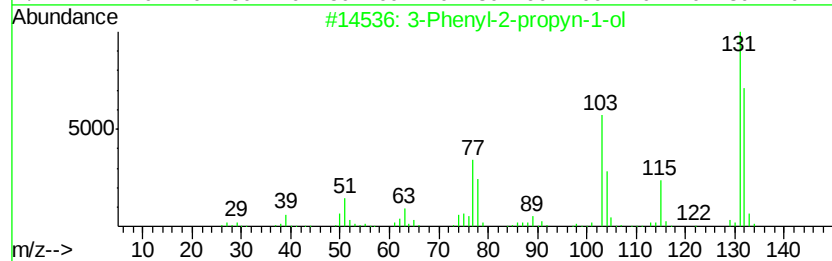
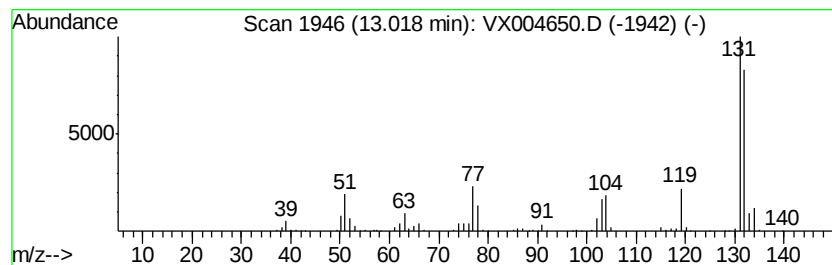
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 2 3-Phenyl-2-propyn-1-ol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.02	6.85 ug/l	134851	1,4-Dichlorobenzene-d4	12.08

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	93
2			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	93
3			Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	93
4			Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	87
5			2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	80



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Operator : JC/MD
Sample : J4898-01 50X
Misc : 5.26g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 7 Sample Multiplier: 1

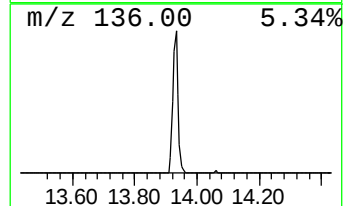
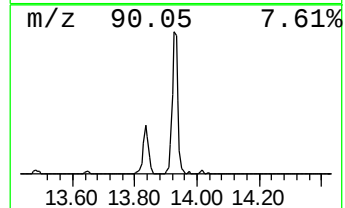
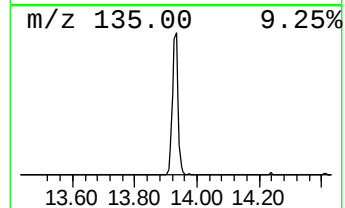
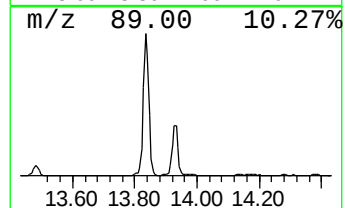
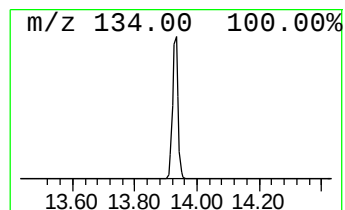
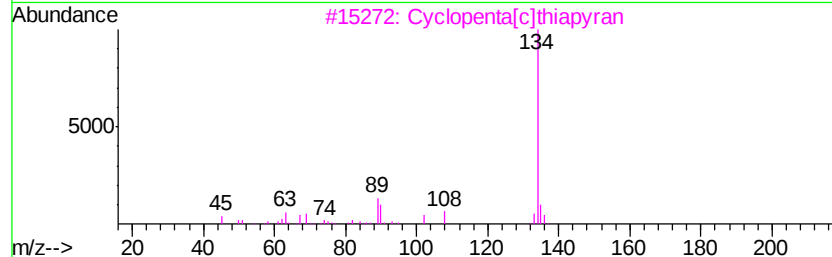
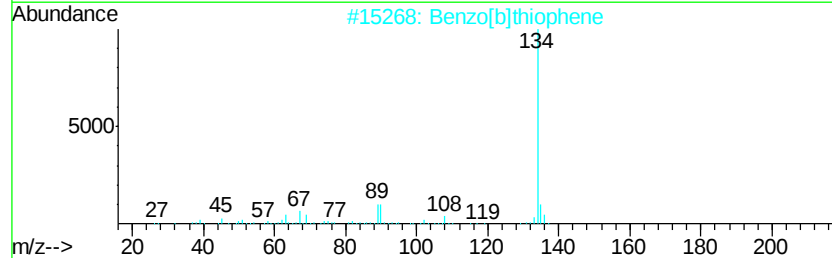
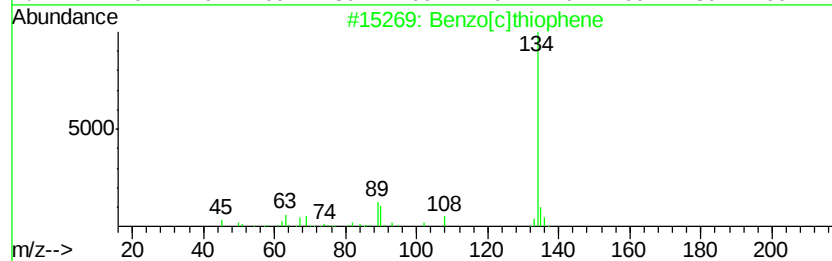
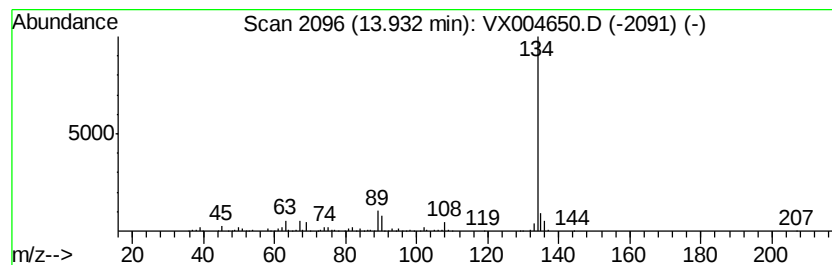
Instrument :
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ClientSampleId :
ROLLOFF01-COMP-FILL

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_X\METHOD\82X091118W.M
Quant Title : SW846 8260

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

Peak Number 3 Benzo[c]thiophene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.93	8.47 ug/l	166900	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	Benzo[c]thiophene	134 C8H6S	000270-82-6	97
2	Benzo[b]thiophene	134 C8H6S	000095-15-8	95
3	Cyclopenta[c]thiapyran	134 C8H6S	000270-63-3	86
4	[1]Benzo[thieno[2',3':3,4]cyclobu...	246 C13H10O3S	034002-19-2	74
5	Thiazolidin-4-one, 5-benzylideno...	287 C13H9N3OS2	351062-07-2	45



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Sample : J4898-01 50X
Misc : 5.26g/5mL/100uL/5.0mL/MSVOA_X/MEOH
ALS Vial : 7 Sample Multiplier: 1

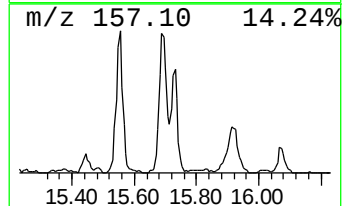
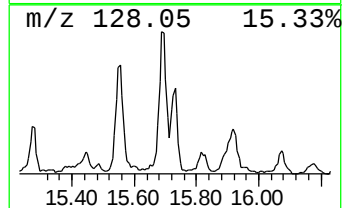
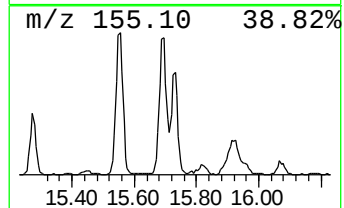
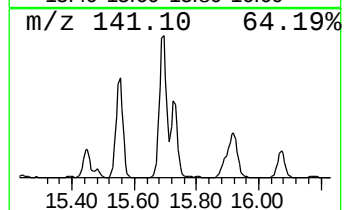
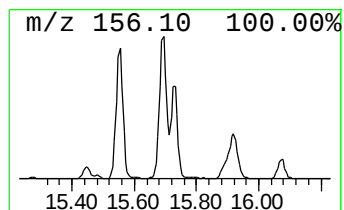
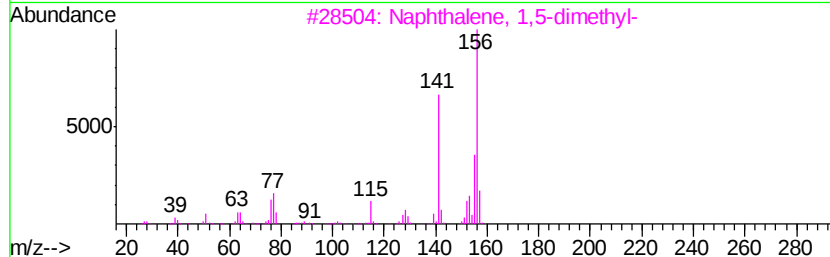
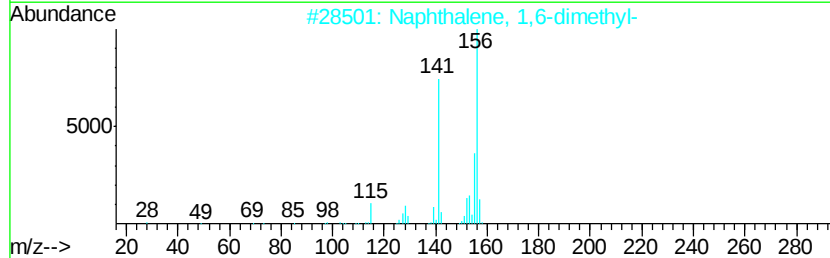
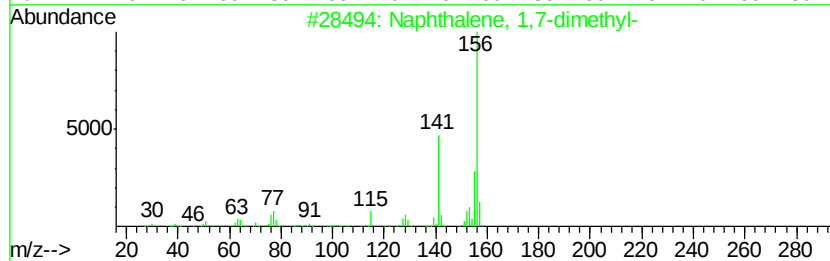
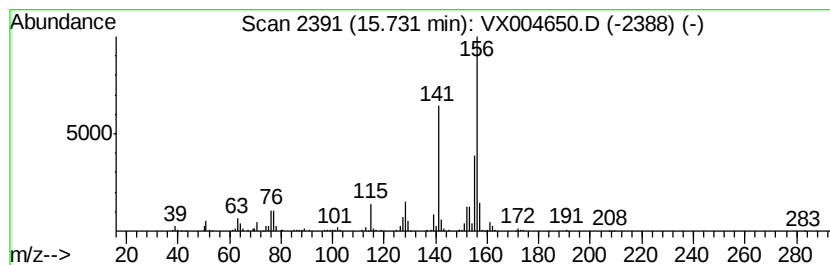
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Peak Number 9 Naphthalene, 1,7-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.73	7.39 ug/l	145511	1,4-Dichlorobenzene-d4	12.08
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,7-dimethyl-	156 C12H12	000575-37-1 96
2		Naphthalene, 1,6-dimethyl-	156 C12H12	000575-43-9 96
3		Naphthalene, 1,5-dimethyl-	156 C12H12	000571-61-9 94
4		Naphthalene, 2,6-dimethyl-	156 C12H12	000581-42-0 94
5		Naphthalene, 2,7-dimethyl-	156 C12H12	000582-16-1 93



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Indene	12.47	38.4	ug/l	756915	4	12.08	985002	50.0
3-Phenyl-2-propyn...	13.02	6.8	ug/l	134851	4	12.08	985002	50.0
Benzo[c]thiophene	13.93	8.5	ug/l	166900	4	12.08	985002	50.0
Naphthalene, 1,7-...	15.73	7.4	ug/l	145511	4	12.08	985002	50.0