

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM092317\
 Data File : BM011710.D
 Acq On : 23 Sep 2017 17:57
 Operator : SJ/JU
 Sample : I5324-05
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AU0

Integration Parameters: LSCINT.P

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

Filtering: 5
 Min Area: 0.1 % 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:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.281	328	339	342	rBV3	64520162	175850894	49.22%	15.120%
2	6.634	563	569	579	rVB	366825	608301	0.17%	0.052%
3	7.104	643	649	659	rVB	214984	377100	0.11%	0.032%
4	7.293	668	681	690	rBV	651512	1281861	0.36%	0.110%
5	7.487	706	714	724	rVB	771090	1394277	0.39%	0.120%
6	7.863	766	778	790	rBV	35101191	89313178	25.00%	7.679%
7	8.069	790	813	817	rBV4	68825403	357259091	100.00%	30.717%
8	8.357	847	862	864	rBV5	65058590	196864706	55.10%	16.926%
9	8.651	905	912	924	rVB	580275	974370	0.27%	0.084%
10	9.122	984	992	994	rBV	932051	1515202	0.42%	0.130%
11	9.175	994	1001	1013	rVB	12349430	21171882	5.93%	1.820%
12	9.457	1042	1049	1058	rBV	800342	1341296	0.38%	0.115%
13	9.769	1097	1102	1108	rVV	420177	771412	0.22%	0.066%
14	9.839	1108	1114	1131	rVB2	833626	1870408	0.52%	0.161%
15	10.375	1198	1205	1214	rBV	1042429	1857077	0.52%	0.160%
16	10.657	1238	1253	1258	rBV2	64353304	165332713	46.28%	14.215%
17	10.769	1266	1272	1279	rBV	995519	1533234	0.43%	0.132%
18	11.157	1327	1338	1346	rBV	28472758	49579755	13.88%	4.263%
19	11.363	1362	1373	1389	rVB	10154323	17349158	4.86%	1.492%
20	11.575	1401	1409	1422	rVB	1747384	3061358	0.86%	0.263%
21	12.045	1481	1489	1502	rVB3	192822	400077	0.11%	0.034%
22	12.739	1597	1607	1609	rBV3	417249	945173	0.26%	0.081%
23	12.780	1609	1614	1622	rVB	1686223	2733561	0.77%	0.235%
24	13.386	1709	1717	1726	rBV	7687884	11649694	3.26%	1.002%
25	13.980	1808	1818	1825	rBV	2185014	3547807	0.99%	0.305%
26	14.274	1862	1868	1875	rVB	2017654	3065988	0.86%	0.264%
27	14.580	1914	1920	1932	rBV2	1179989	1985848	0.56%	0.171%
28	15.574	2082	2089	2095	rBV	2536047	3861893	1.08%	0.332%
29	15.686	2102	2108	2120	rVB	1115682	1583248	0.44%	0.136%
30	17.321	2378	2386	2396	rBV2	1517738	2246919	0.63%	0.193%
31	17.421	2396	2403	2414	rBV	3134241	4300846	1.20%	0.370%
32	18.692	2608	2619	2627	rBV	12857880	17463754	4.89%	1.502%
33	18.815	2635	2640	2650	rVB	722498	1011251	0.28%	0.087%
34	18.980	2663	2668	2671	rBV2	373953	657239	0.18%	0.057%

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Integration Parameters: LSCINT.P

Integrator: RTE
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Max Peaks: 100
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If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.M
Title : SVOA CALIBRATION

35	19.021	2671	2675	2679	rVV	842721	1328607	0.37%	0.114%
36	19.068	2679	2683	2688	rVB2	326192	523820	0.15%	0.045%
37	19.709	2786	2792	2802	rBV	3653214	5206340	1.46%	0.448%
38	21.486	3089	3094	3100	rBV	2093476	2624104	0.73%	0.226%
39	23.697	3463	3470	3482	rVB2	2778701	5626790	1.57%	0.484%
40	23.850	3490	3496	3505	rVB	1248146	2591640	0.73%	0.223%
41	24.080	3531	3535	3541	rVB	216818	398466	0.11%	0.034%

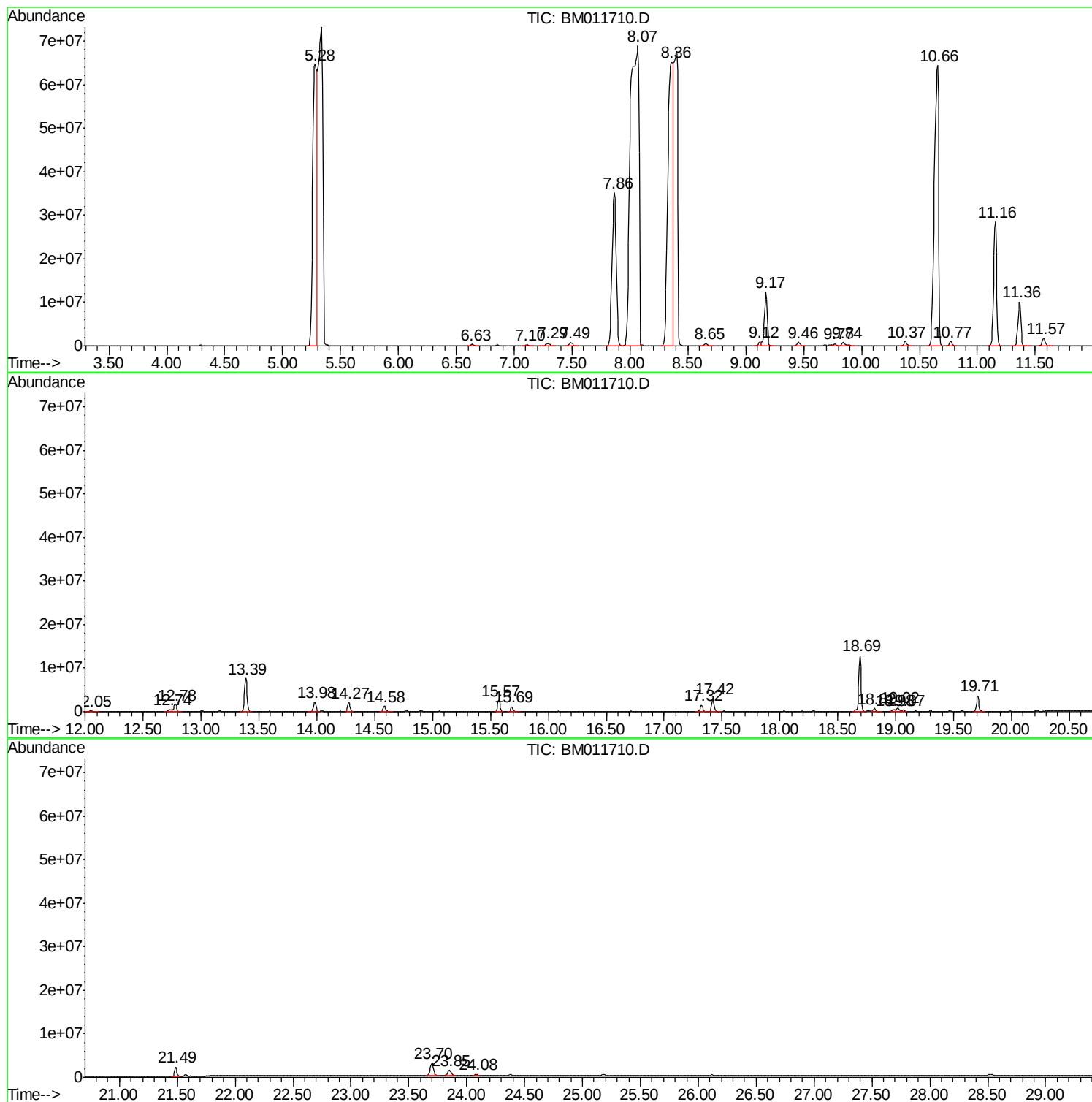
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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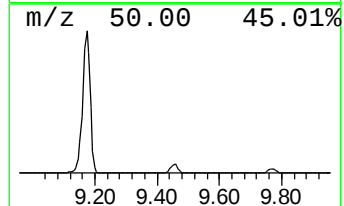
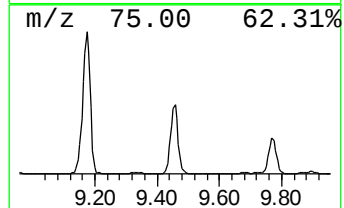
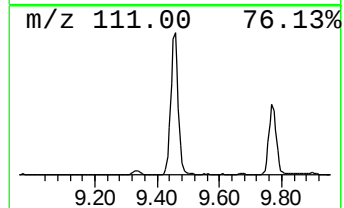
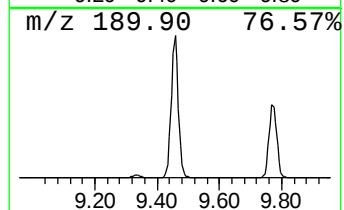
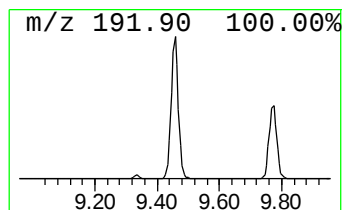
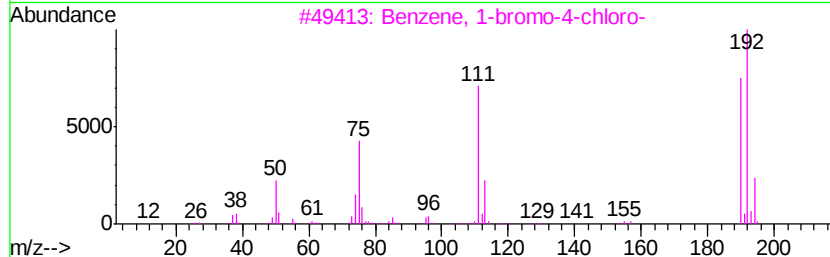
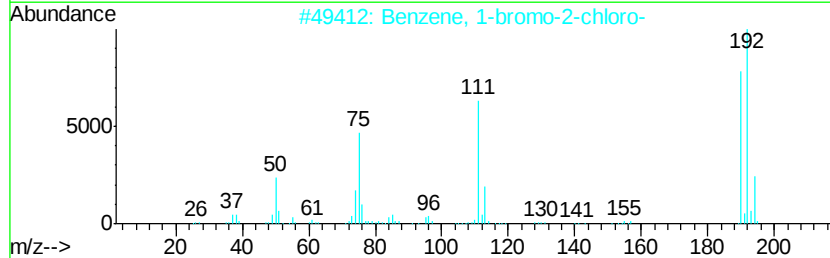
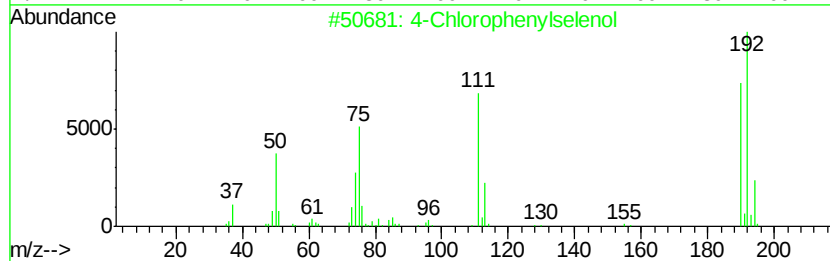
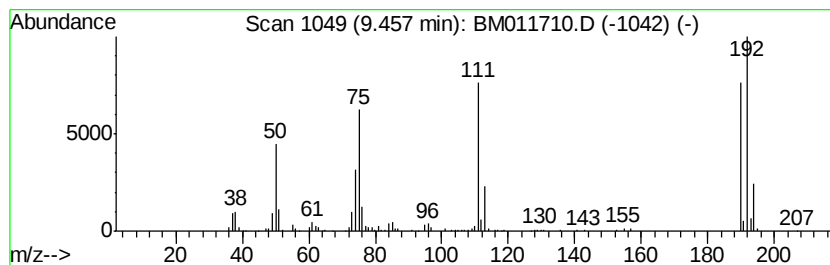
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 4-Chlorophenylselenol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.46	17.50 ng/ul	1341300	Napthalene-d8	10.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	99
2		Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
3		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	97
4		Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	93
5		Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	93



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Instrument :
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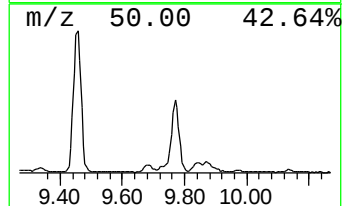
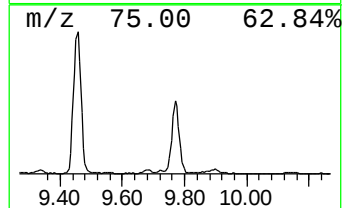
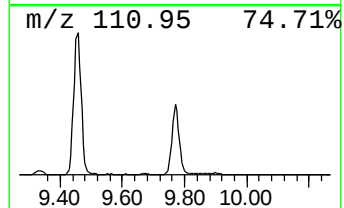
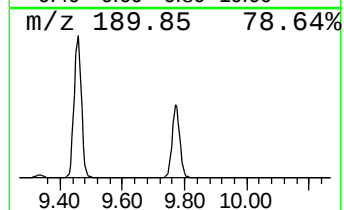
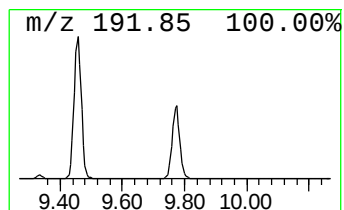
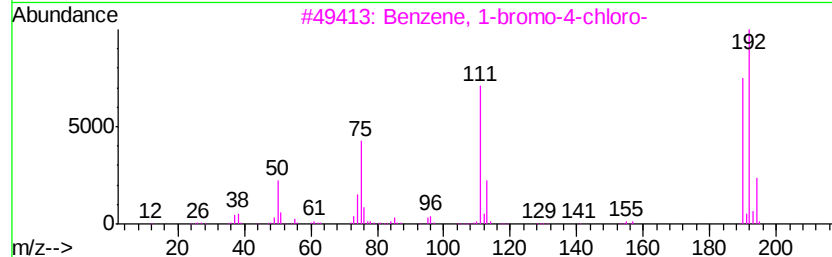
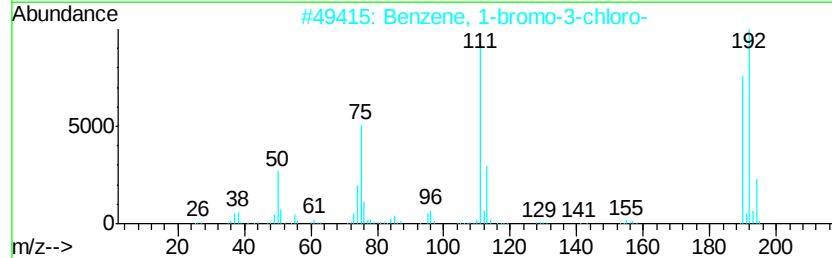
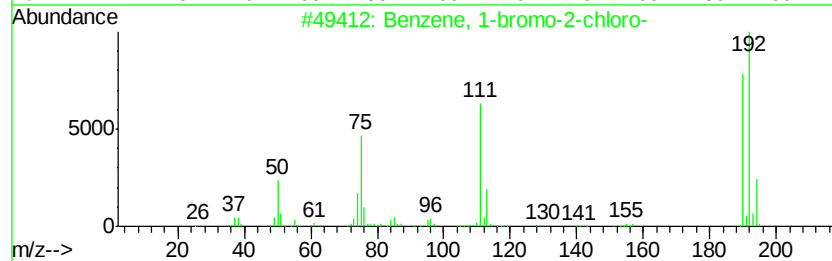
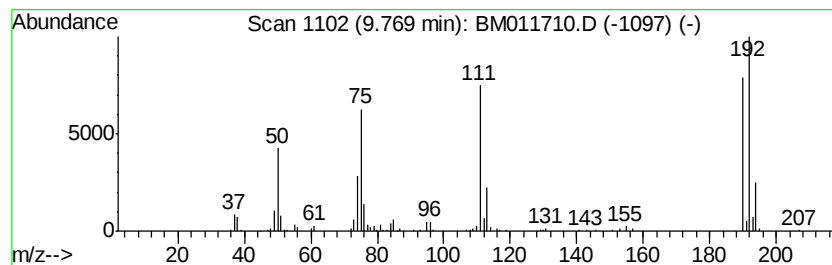
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.M
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Benzene, 1-bromo-2-chloro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.77	10.06 ng/ul	771412	Napthalene-d8	10.77

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
2			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	98
3			Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
4			Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	98
5			Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	97



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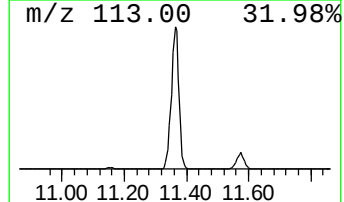
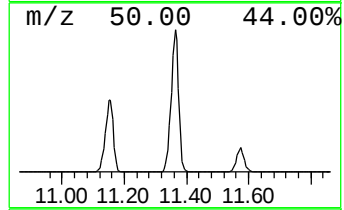
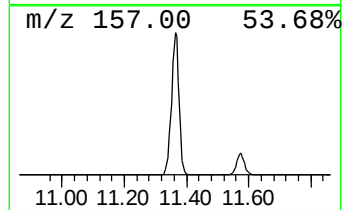
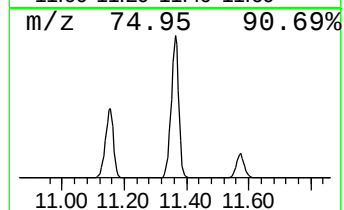
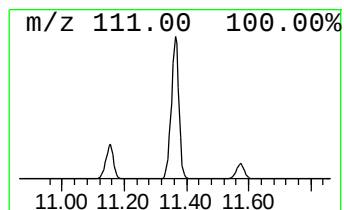
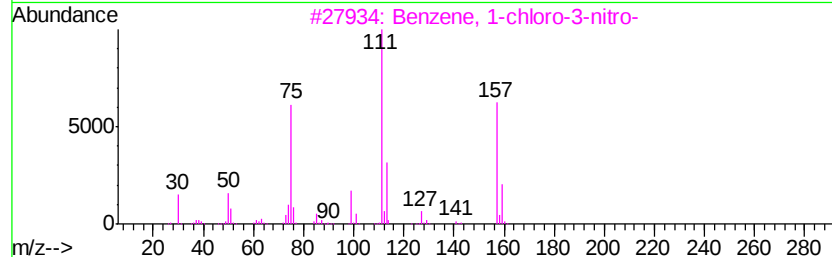
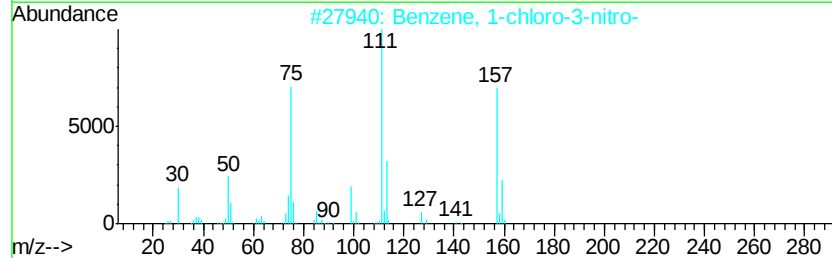
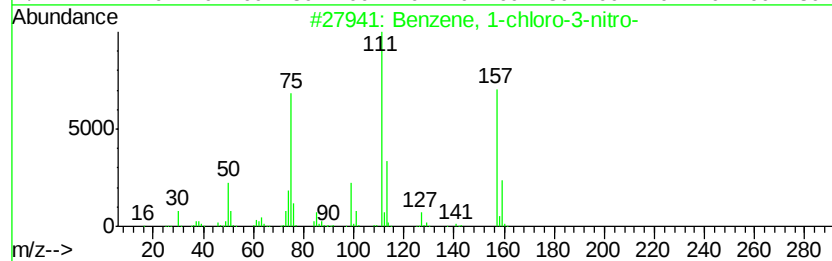
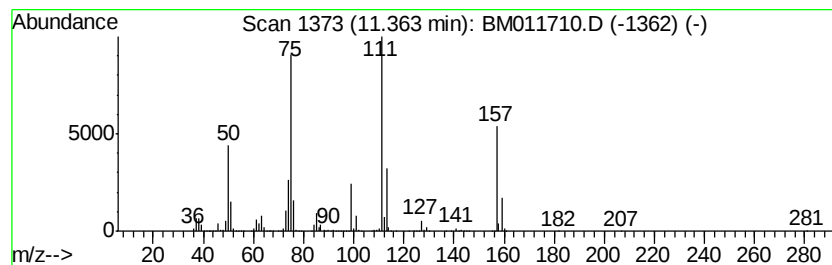
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Benzene, 1-chloro-3-nitro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.36	226.31 ng/ul	17349200	Naphthalene-d8	10.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	97
2		Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	91
3		Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	90
4		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	64
5		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	58



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Instrument :
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ClientSampled :
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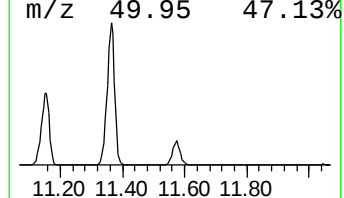
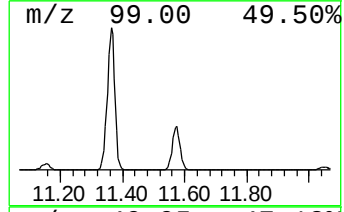
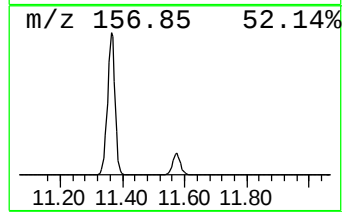
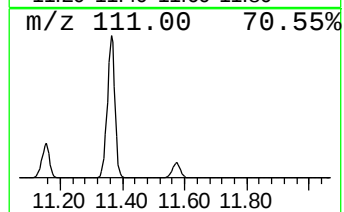
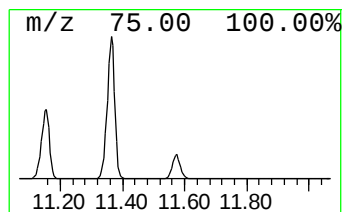
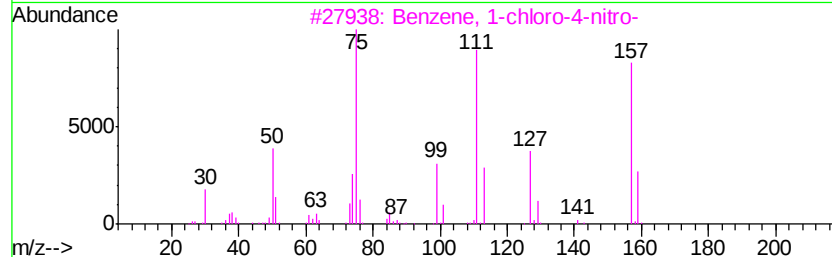
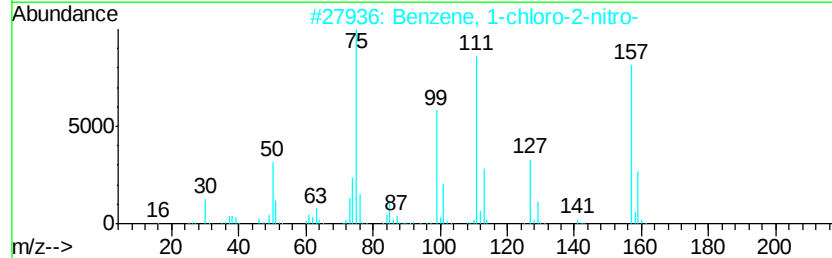
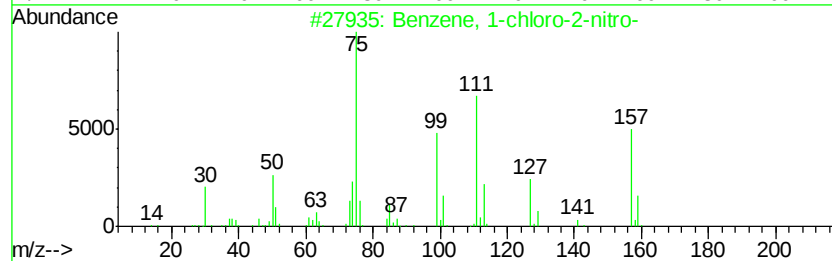
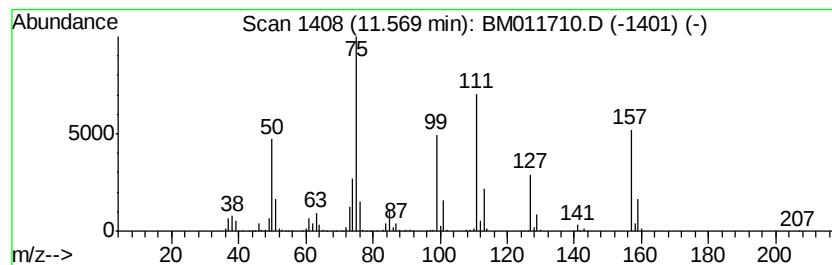
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TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Benzene, 1-chloro-2-nitro- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	39.93 ng/ul	3061360	Naphthalene-d8	10.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	98
2		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	97
3		Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	97
4		Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	93
5		Benzene, 1-chloro-2-nitro-	157	C6H4ClNO2	000088-73-3	90



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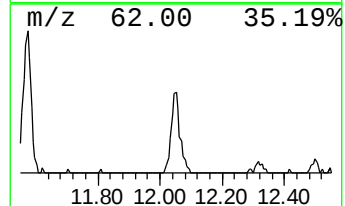
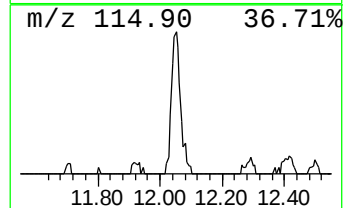
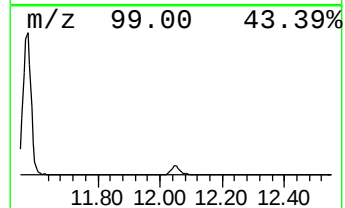
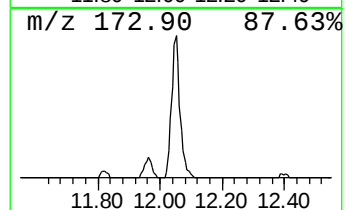
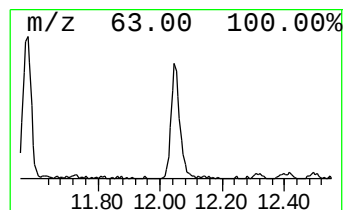
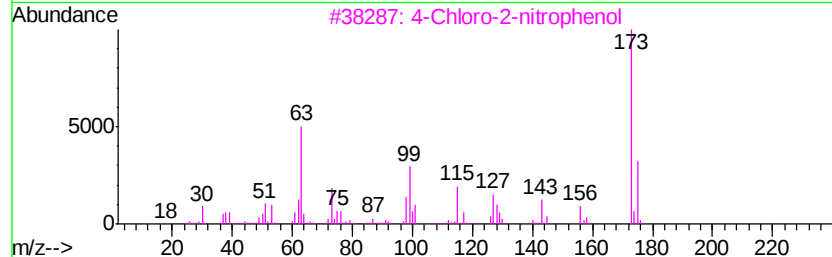
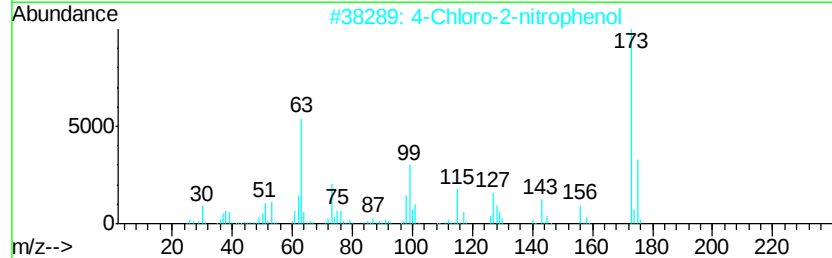
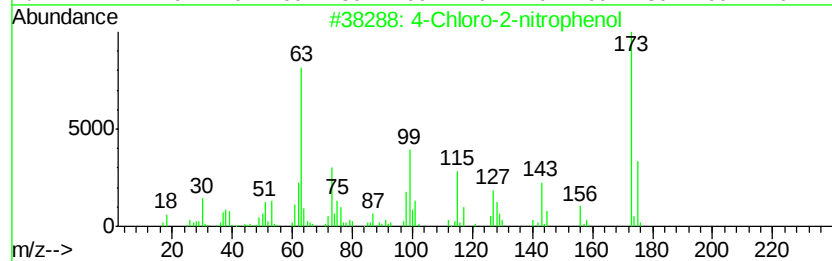
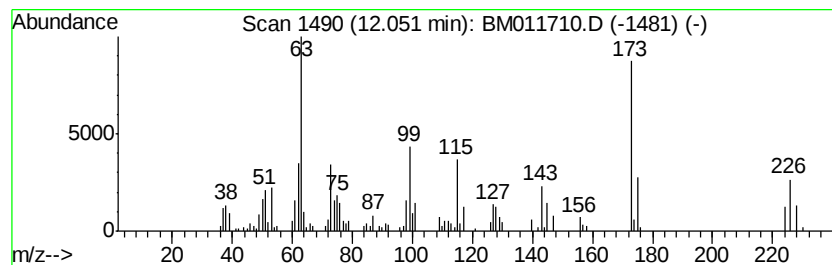
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 4-Chloro-2-nitrophenol Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.05	5.22 ng/ul	400077	Naphthalene-d8	10.77

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	99
2		4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	95
3		4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	95
4		Phenol, 2-chloro-4-nitro-	173	C6H4ClNO3	000619-08-9	80
5		Phenol, 5-chloro-2-nitro-	173	C6H4ClNO3	000611-07-4	55



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM092317\
Data File : BM011710.D
Acq On : 23 Sep 2017 17:57
Operator : SJ/JU
Sample : I5324-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

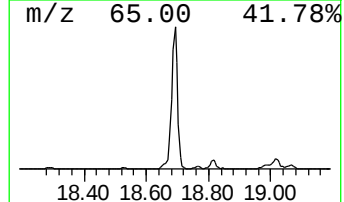
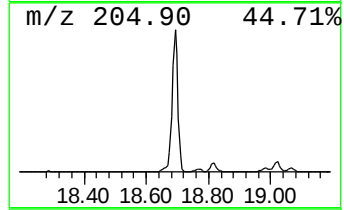
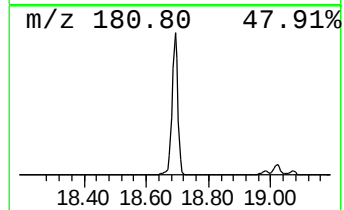
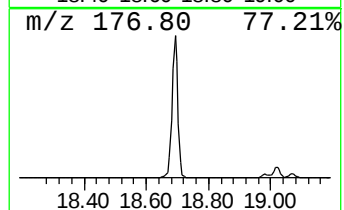
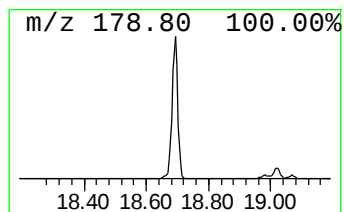
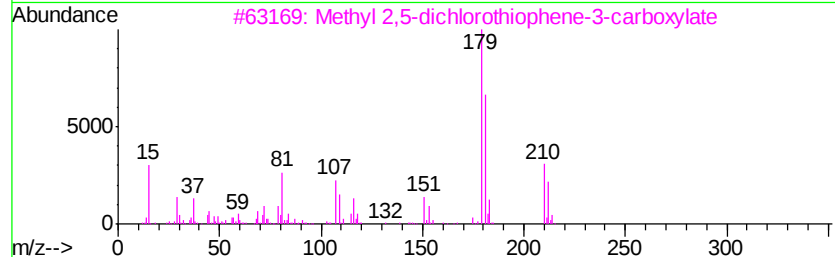
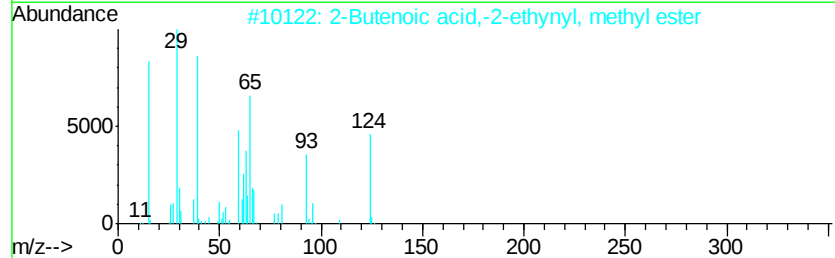
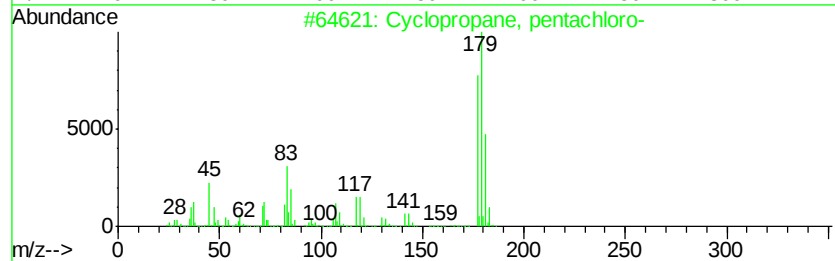
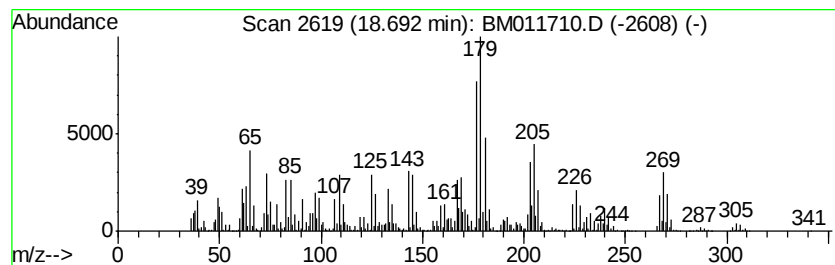
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.69	155.45 ng/ul	17463800	Phenanthrene-d10	17.32		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopropane, pentachloro-	212	C3HC15	006262-51-7	22
2		2-Butenoic acid, -2-ethynyl, meth...	124	C7H8O2	1000222-03-1	17
3		Methyl 2,5-dichlorothiophene-3-c...	210	C6H4Cl2O2S	145129-54-0	14
4		4-Chloro-3-quinolinol	179	C9H6ClNO	032435-60-2	11
5		Cloxyquin	179	C9H6ClNO	000130-16-5	11



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM092317\
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Operator : SJ/JU
Sample : I5324-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

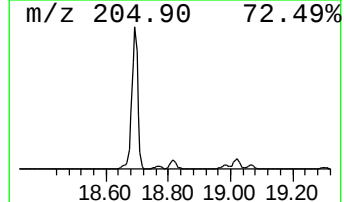
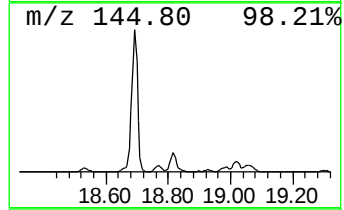
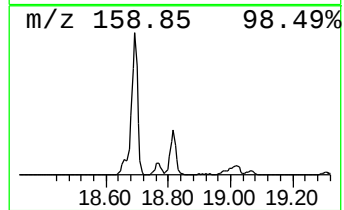
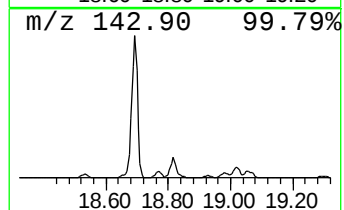
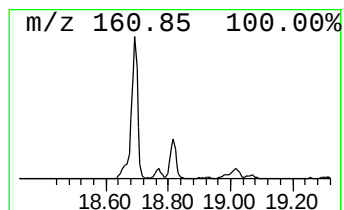
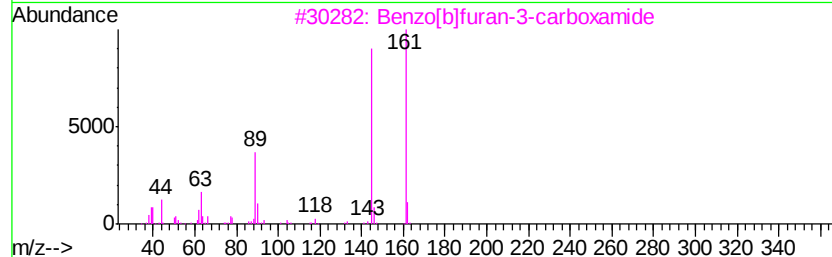
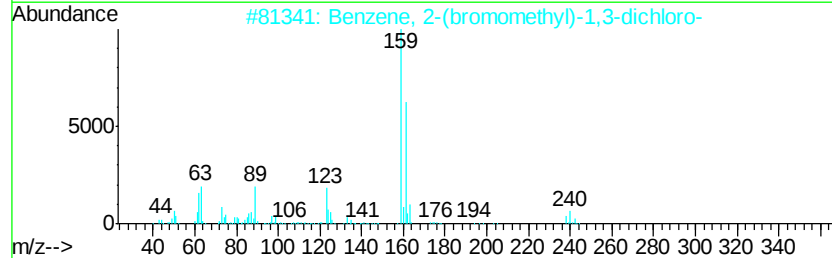
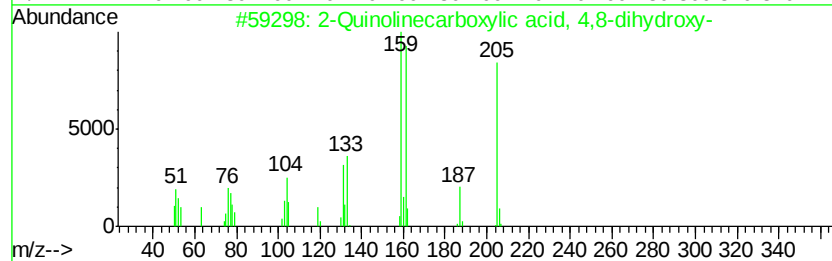
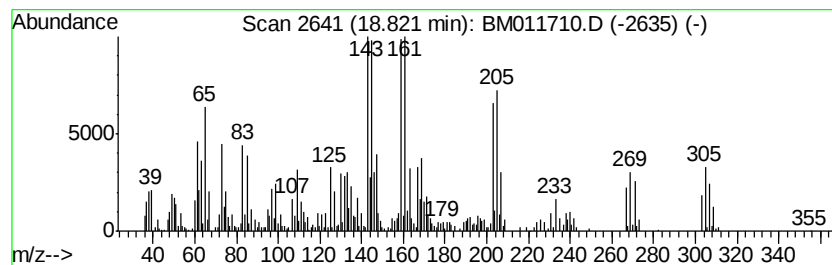
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown-02 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.82	9.00 ng/ul	1011250	Phenanthrene-d10	17.32	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Quinolinecarboxylic acid, 4,8-...	205	C10H7N04	000059-00-7	35
2	Benzene, 2-(bromomethyl)-1,3-dic...	238	C7H5BrCl2	020443-98-5	25
3	Benzo[b]furan-3-carboxamide	161	C9H7N02	1000262-22-7	11
4	Benzene, 1-methyl-2-nitroso-	121	C7H7NO	000611-23-4	11
5	Aniline, 3,5-bis(3-tolyloxy)-	305	C20H19N02	329726-24-1	10



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM092317\
Data File : BM011710.D
Acq On : 23 Sep 2017 17:57
Operator : SJ/JU
Sample : I5324-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

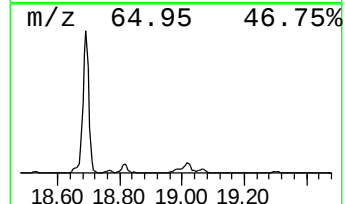
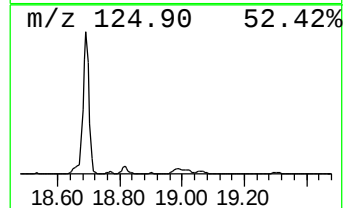
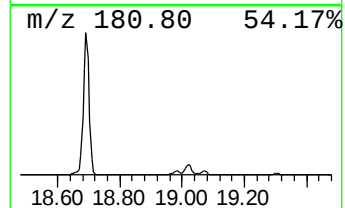
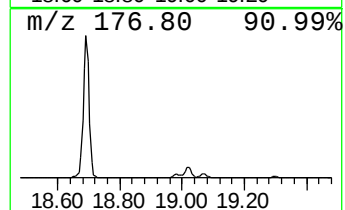
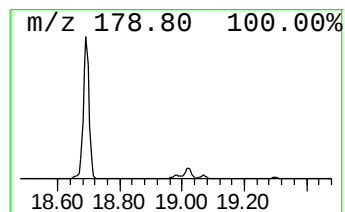
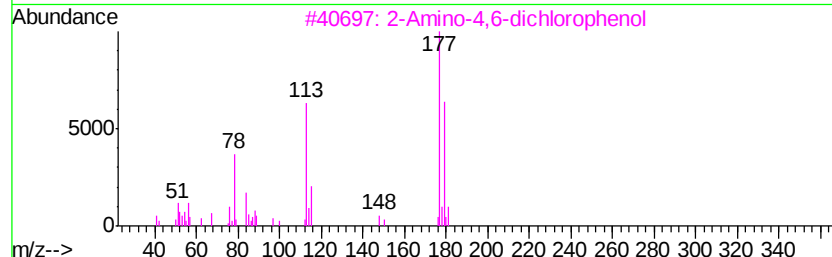
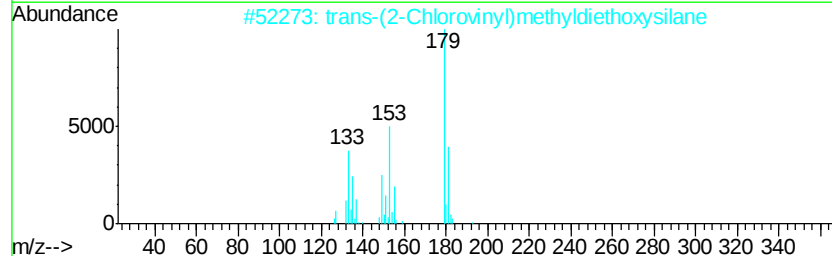
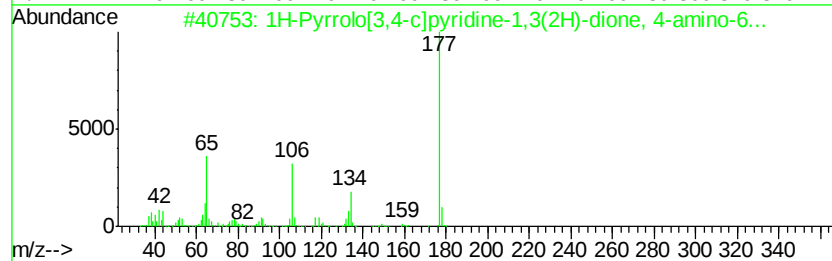
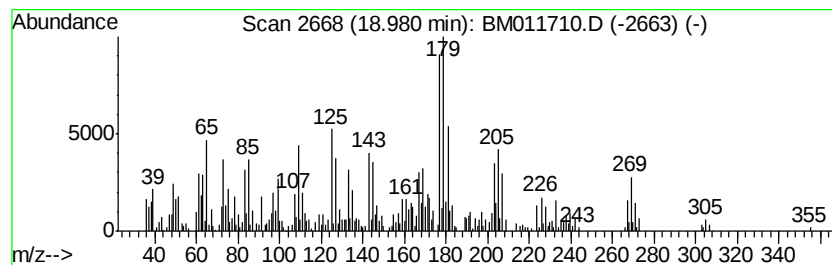
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Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 unknown-03 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.		
18.98	5.85 ng/ul	657239	Phenanthrene-d10	17.32		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Pyrrolo[3,4-c]pyridine-1,3(2H...	177	C8H7N3O2	090004-20-9	22
2		trans-(2-Chlorovinyl)methyldieth...	194	C7H15ClO2Si	1000139-96-9	16
3		2-Amino-4,6-dichlorophenol	177	C6H5Cl2NO	000527-62-8	11
4		2(3H)-Benzothiazolone, 3-methyl-...	179	C8H9N3S	001128-67-2	11
5		Methyl 2,5-dichlorothiophene-3-c...	210	C6H4Cl2O2S	145129-54-0	10



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM092317\
Data File : BM011710.D
Acq On : 23 Sep 2017 17:57
Operator : SJ/JU
Sample : I5324-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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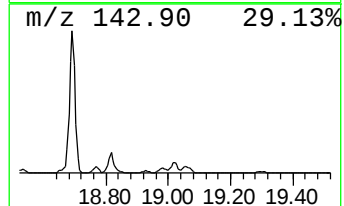
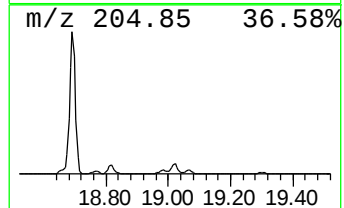
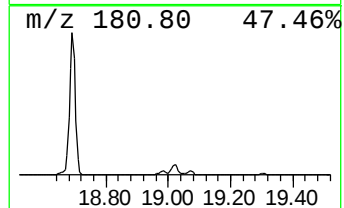
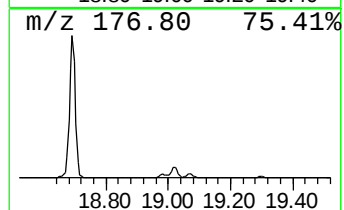
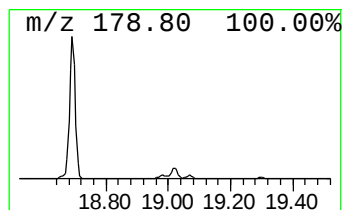
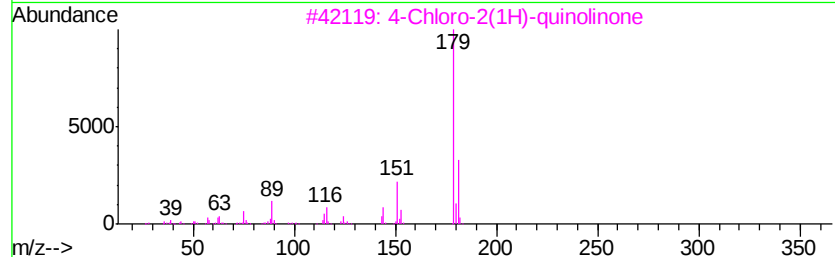
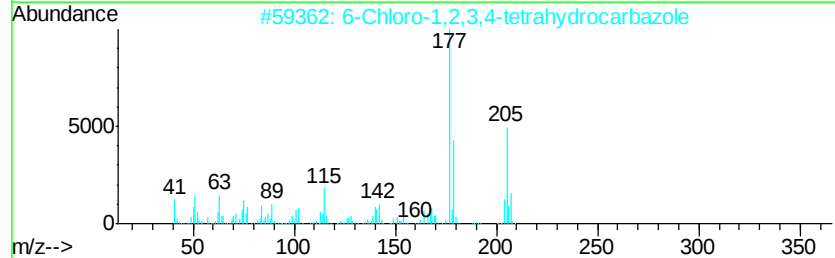
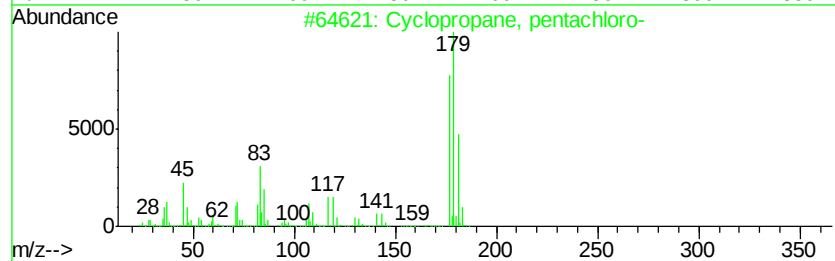
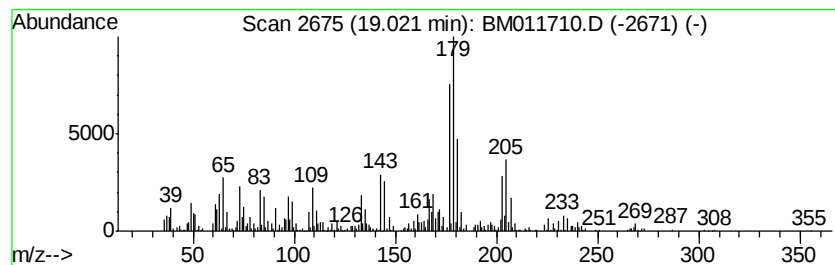
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 unknown-04 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.02	11.83 ng/ul	1328610	Phenanthrene-d10	17.32

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Cyclopropane, pentachloro-	212	C3HCl5	006262-51-7	37
2		6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	25
3		4-Chloro-2(1H)-quinolinone	179	C9H6ClNO	020146-59-2	22
4		Octamethyl-1,4-cyclohexadiene	192	C14H24	172684-93-4	15
5		3-Buten-2-one, 4-(2,6,6-trimethy...	192	C13H20O	014901-07-6	15



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM092317\
Data File : BM011710.D
Acq On : 23 Sep 2017 17:57
Operator : SJ/JU
Sample : I5324-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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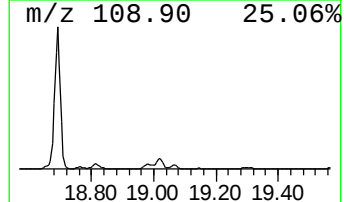
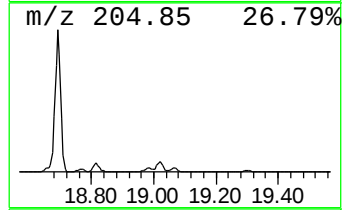
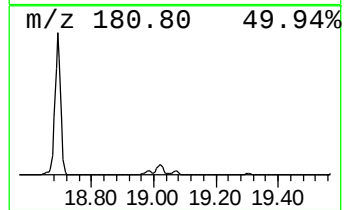
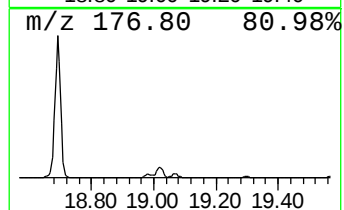
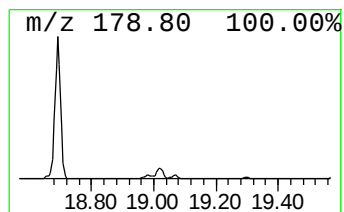
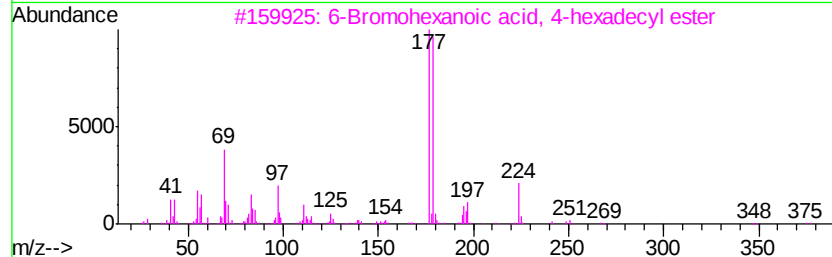
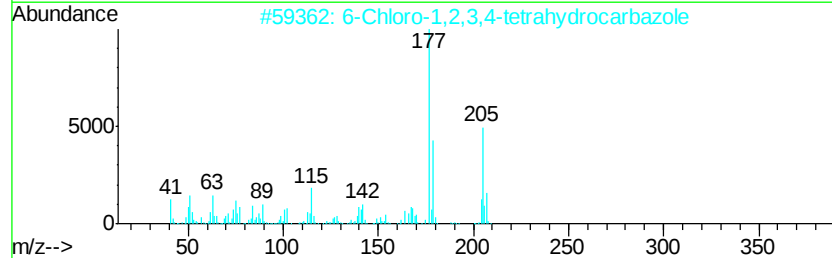
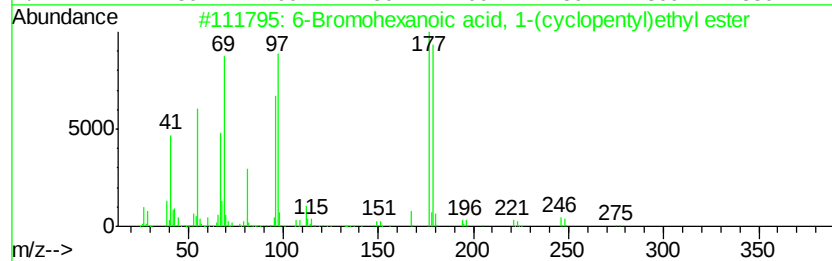
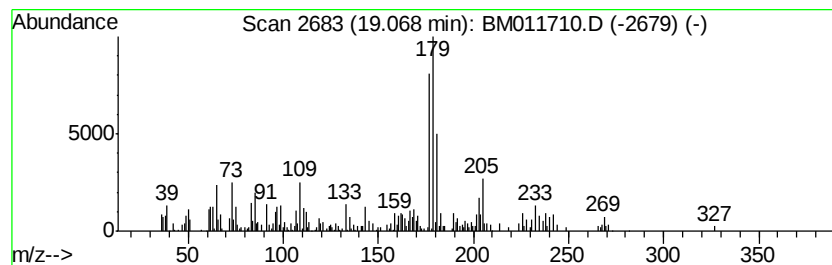
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 unknown-05 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.07	4.66 ng/ul	523820	Phenanthrene-d10	17.32

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			6-Bromohexanoic acid, 1-(cyclope...	290	C13H23BrO2	1000293-29-6	25
2			6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	25
3			6-Bromohexanoic acid, 4-hexadecy...	418	C22H43BrO2	1000282-90-6	22
4			Silane, trimethyl(1-phenylethoxy)-	194	C11H18OSi	014856-75-8	18
5			6-Chloro-2-methylquinoline	177	C10H8ClN	000092-46-6	14



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM092317\
Data File : BM011710.D
Acq On : 23 Sep 2017 17:57
Operator : SJ/JU
Sample : I5324-05
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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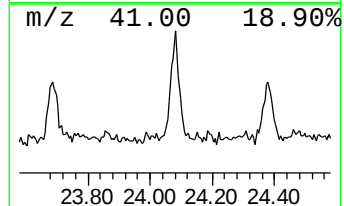
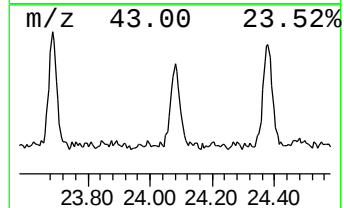
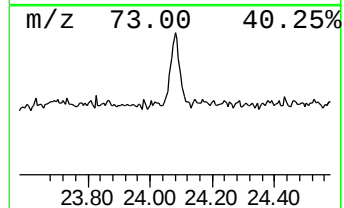
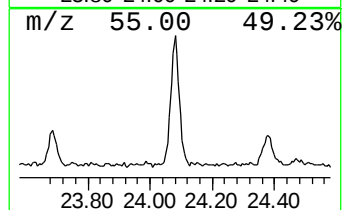
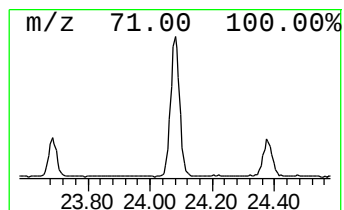
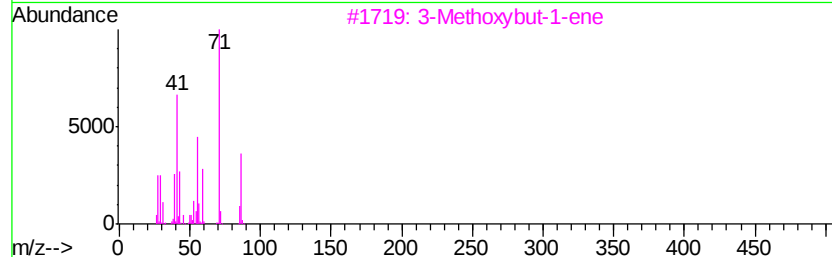
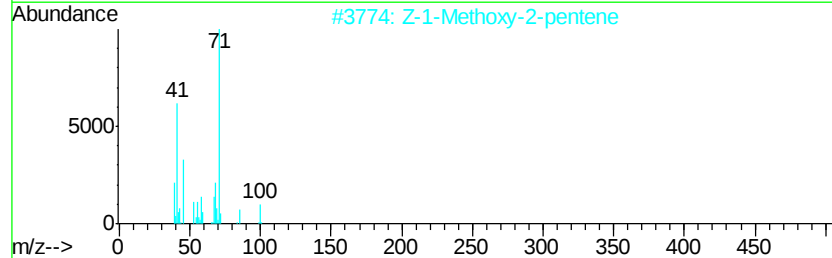
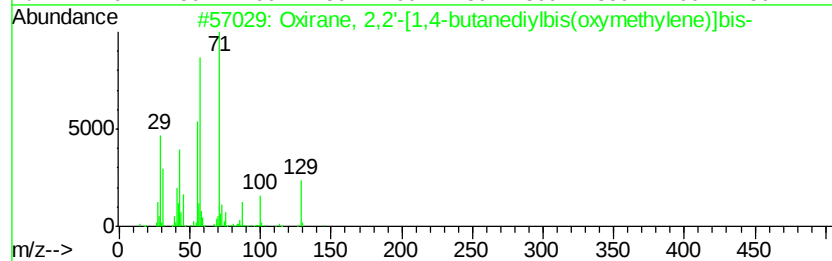
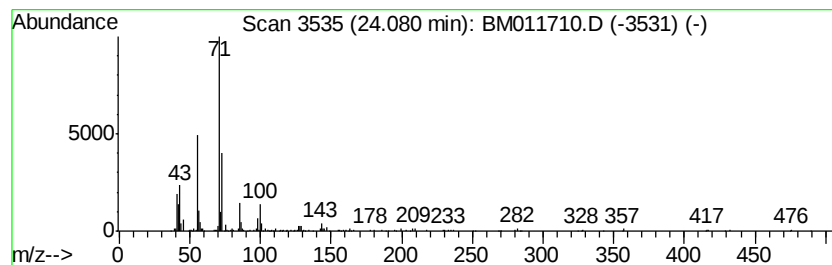
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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 Oxirane, 2,2'-[1,4-butanedi... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.08	3.08 ng/ul	398466	Perylene-d12	23.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Oxirane, 2,2'-[1,4-butanediylbis...	202	C10H18O4	002425-79-8	56
2		Z-1-Methoxy-2-pentene	100	C6H12O	1000279-59-4	50
3		3-Methoxybut-1-ene	86	C5H10O	017351-24-5	45
4		Cis-3-methylpent-3-ene-5-ol	100	C6H12O	030804-75-2	45
5		Butanoic acid, 2-butoxy-1-methyl...	216	C11H20O4	007492-70-8	42



Data Path : Z:\HPCHEM1\BNA_M\DATA\BM092317\
 Data File : BM011710.D
 Acq On : 23 Sep 2017 17:57
 Operator : SJ/JU
 Sample : I5324-05
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 C0AU0

Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM090817MA.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
4-Chlorophenylsel...	9.46	17.5	ng/ul	1341300	2	10.77	1533230	20.0
Benzene, 1-bromo-...	9.77	10.1	ng/ul	771412	2	10.77	1533230	20.0
Benzene, 1-chloro...	11.36	226.3	ng/ul	17349200	2	10.77	1533230	20.0
Benzene, 1-chloro...	11.57	39.9	ng/ul	3061360	2	10.77	1533230	20.0
4-Chloro-2-nitrop...	12.05	5.2	ng/ul	400077	2	10.77	1533230	20.0
unknown-01	18.69	155.4	ng/ul	17463800	4	17.32	2246920	20.0
unknown-02	18.82	9.0	ng/ul	1011250	4	17.32	2246920	20.0
unknown-03	18.98	5.8	ng/ul	657239	4	17.32	2246920	20.0
unknown-04	19.02	11.8	ng/ul	1328610	4	17.32	2246920	20.0
unknown-05	19.07	4.7	ng/ul	523820	4	17.32	2246920	20.0
Oxirane, 2,2'-[1,...	24.08	3.1	ng/ul	398466	6	23.85	2591640	20.0