LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092118\
Data File : BM016844.D
Acq On : 21 Sep 2018 22:18
Operator : SJ/JU
Sample : J5013-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08R4

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.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	3.016	3	5	11	rVB	2416162	2762269	1.30%	0.254%
2	5.093	348	358	361	rBV4	50313790	127273152	60.07%	11.705%
3	6.422	575	584	592	rVB	617703	1000021	0.47%	0.092%
4	6.675	614	627	631	rBV	208666	424928	0.20%	0.039%
5	6.716	631	634	641	rVB	144226	227467	0.11%	0.021%
6	6.898	658	665	677	rVB	266184	491034	0.23%	0.045%
7	7.092	687	698	708	rBV	541807	1575155	0.74%	0.145%
8	7.269	720	728	742	rVB	771945	1726655	0.81%	0.159%
9	7.463	751	761	772	rVB3	80046	276047	0.13%	0.025%
10	7.639	779	791	805	rBV	29214526	127123025	60.00%	11.692%
11	7.798	805	818	820	rBV4	53341522	161313691	76.14%	14.836%
12	8.139	861	876	878	rBV4	53809287	186724091	88.13%	17.173%
13	8.434	920	926	934	rVB	865064	1346388	0.64%	0.124%
14	8.886	996	1003	1006	rVV	1090180	2027562	0.96%	0.186%
15	8.945	1006	1013	1027	rVB	13721579	24804056	11.71%	2.281%
16	9.216	1051	1059	1070	rBV	1523138	2408444	1.14%	0.222%
17	9.439	1089	1097	1101	rBV	156927	261751	0.12%	0.024%
18	9.528	1107	1112	1117	rVB	695150	1185347	0.56%	0.109%
19	9.598	1117	1124	1132	rBV2	964089	2012847	0.95%	0.185%
20	10.128	1207	1214	1222	rBV	1408727	2504372	1.18%	0.230%
21	10.416	1248	1263	1266	rBV6	58000045	211872107	100.00%	19.486%
22	10.528	1276	1282	1288	rBV	1558437	2320697	1.10%	0.213%
23	10.916	1335	1348	1354	rBV	46797991	103679760	48.94%	9.536%
24	11.110	1372	1381	1391	rBV	10209643	17729864	8.37%	1.631%
25	11.322	1407	1417	1425	rBV	2068841	3582900	1.69%	0.330%
26	11.798	1487	1498	1507	rVB2	462949	869438	0.41%	0.080%
27	12.486	1604	1615	1619	rBV4	540186	1223023	0.58%	0.112%
28	12.539	1619	1624	1632	rVB	2657812	4275308	2.02%	0.393%
29	12.757	1655	1661	1673	rVB3	121786	286766	0.14%	0.026%
30	12.927	1684	1690	1696	rBV	263198	401429	0.19%	0.037%
31	13.151	1721	1728	1738	rVB	10648130	16352431	7.72%	1.504%
32	13.363	1760	1764	1770	rVB	247482	351122	0.17%	0.032%
33	13.745	1822	1829	1830	rBV	663130	904428	0.43%	0.083%
34	13.774	1830	1834	1839	rVV	2758707	3783399	1.79%	0.348%

LSC Area Percent Report

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Data File : BM016844.D
Acq On : 21 Sep 2018 22:18
Operator : SJ/JU
Sample : J5013-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08R4

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Title : SVOA CALIBRATION

35	14.051	1875	1881	1889	rVV	3333036	4808707	2.27%	0.442%
36	14.357	1928	1933	1945	rVB2	1885091	2983258	1.41%	0.274%
37	14.551	1962	1966	1978	rVV3	294340	465271	0.22%	0.043%
38	14.674	1978	1987	1993	rVB	139718	213293	0.10%	0.020%
39	15.351	2095	2102	2109	rBV	4089365	5948371	2.81%	0.547%
40	15.468	2116	2122	2132	rBV	1111718	1541403	0.73%	0.142%
41	16.504	2294	2298	2303	rVB2	215188	288843	0.14%	0.027%
42	17.104	2391	2400	2406	rBV2	2302285	3331705	1.57%	0.306%
43	17.204	2411	2417	2428	rVV	4503684	6181704	2.92%	0.569%
44	17.298	2428	2433	2437	rVV2	193109	352958	0.17%	0.032%
45	17.339	2437	2440	2448	rVB2	208779	360838	0.17%	0.033%
46	17.821	2513	2522	2527	rVB3	201588	323979	0.15%	0.030%
47	18.074	2559	2565	2571	rVV4	174214	287634	0.14%	0.026%
48	18.480	2629	2634	2643	rVB	12502452	17294125	8.16%	1.591%
49	18.556	2643	2647	2651	rVV3	176810	242208	0.11%	0.022%
50	18.603	2651	2655	2665	rVB	634311	1002827	0.47%	0.092%
51	18.780	2679	2685	2686	rBV2	348959	525562	0.25%	0.048%
52	18.809	2686	2690	2695	rVV	1022861	1646141	0.78%	0.151%
53	18.856	2695	2698	2704	rVB2	260510	387557	0.18%	0.036%
54	19.362	2780	2784	2794	rVB2	179297	292654	0.14%	0.027%
55	19.498	2801	2807	2818	rBV	5333493	7475402	3.53%	0.688%
56	21.292	3107	3112	3120	rBV	3263634	4095073	1.93%	0.377%
57	23.386	3461	3468	3475	rBV	4245255	7949700	3.75%	0.731%
58	23.527	3486	3492	3503	rVB	2307636	4201709	1.98%	0.386%

Sum of corrected areas: 1087301896

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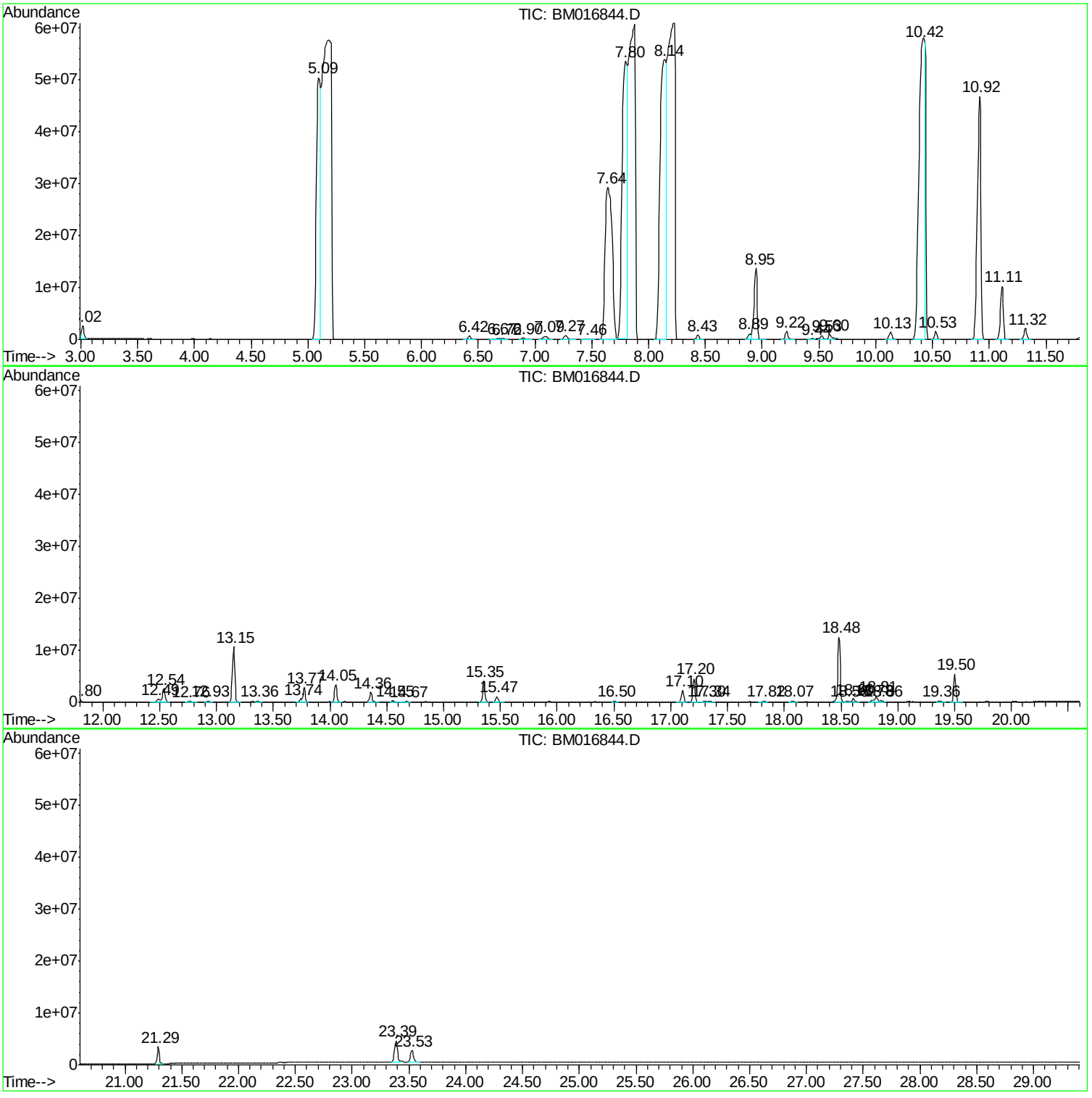
Sample : J5013-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C08R4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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



Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092118\
Data File : BM016844.D
Acq On : 21 Sep 2018 22:18
Operator : SJ/JU
Sample : J5013-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08R4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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

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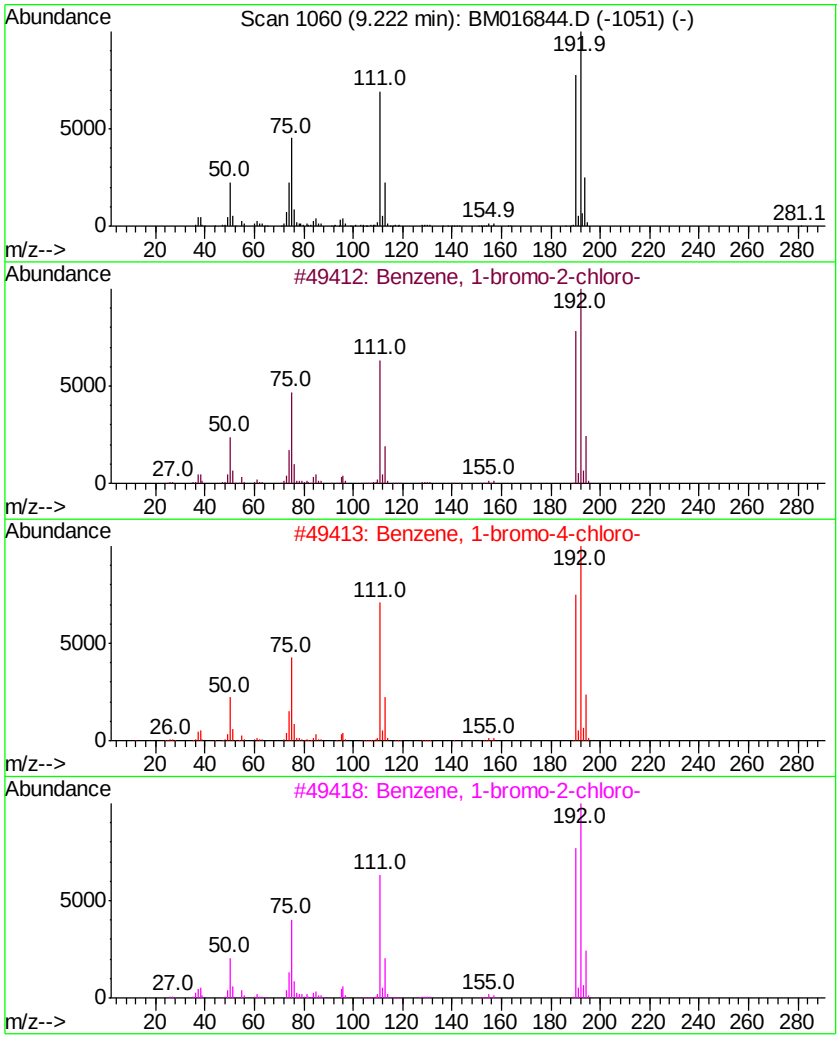
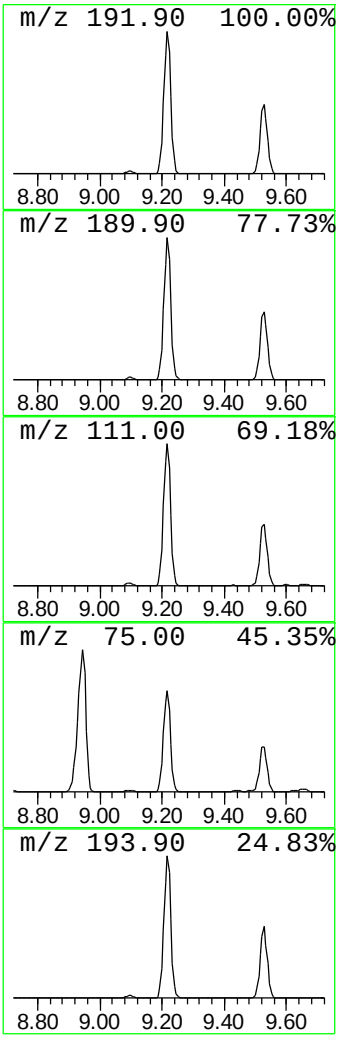
Peak Number 4 Benzene, 1-bromo-2-chloro- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.22	20.76 ng/ul	2408440	Naphthalene-d8	10.53

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





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092118\
Data File : BM016844.D
Acq On : 21 Sep 2018 22:18
Operator : SJ/JU
Sample : J5013-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08R4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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

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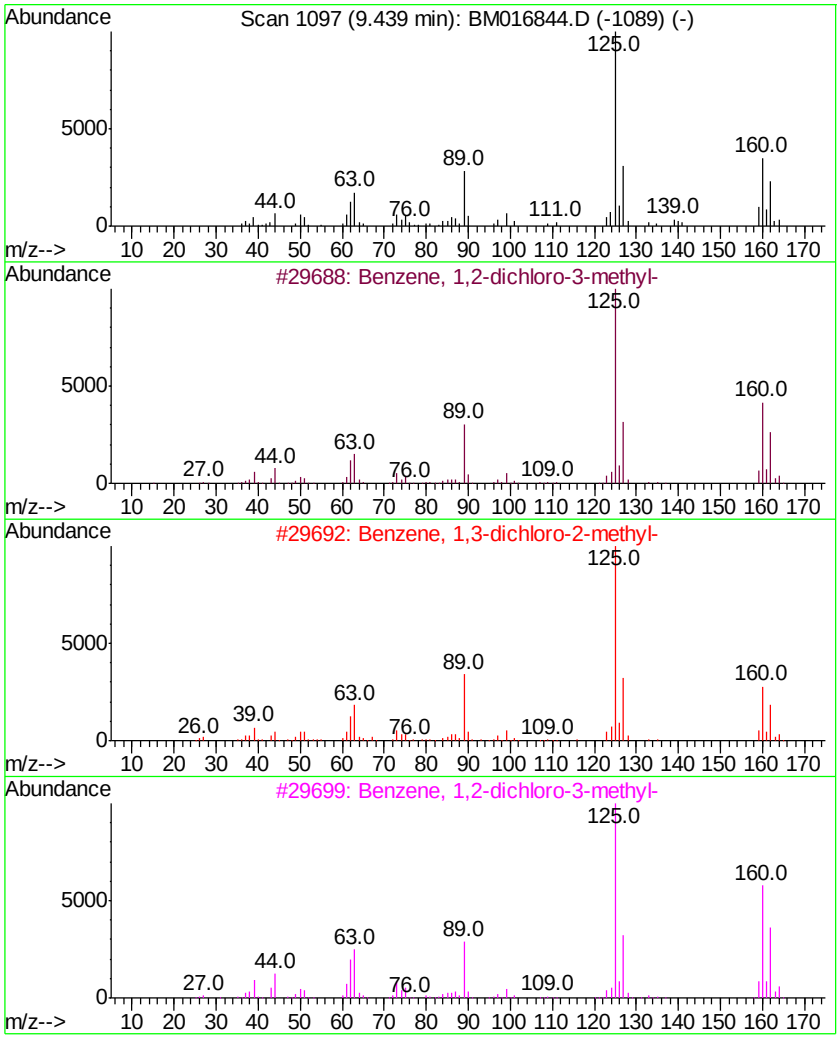
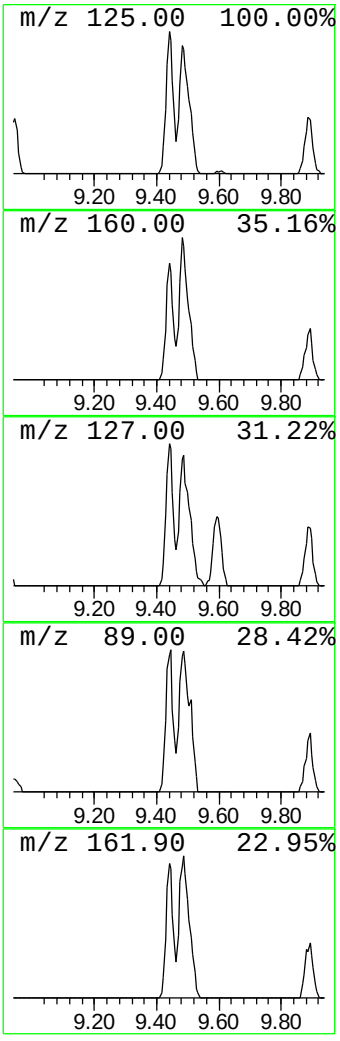
Peak Number 5 Benzene, 1,2-dichloro-3-met... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.44	2.26 ng/ul	261751	Naphthalene-d8	10.53

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	98
2	Benzene, 1,3-dichloro-2-methyl-	160	C7H6Cl2	000118-69-4	97
3	Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	97
4	Benzene, 1,3-dichloro-2-methyl-	160	C7H6Cl2	000118-69-4	97
5	Benzene, 1,3-dichloro-2-methyl-	160	C7H6Cl2	000118-69-4	97





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092118\
Data File : BM016844.D
Acq On : 21 Sep 2018 22:18
Operator : SJ/JU
Sample : J5013-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08R4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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

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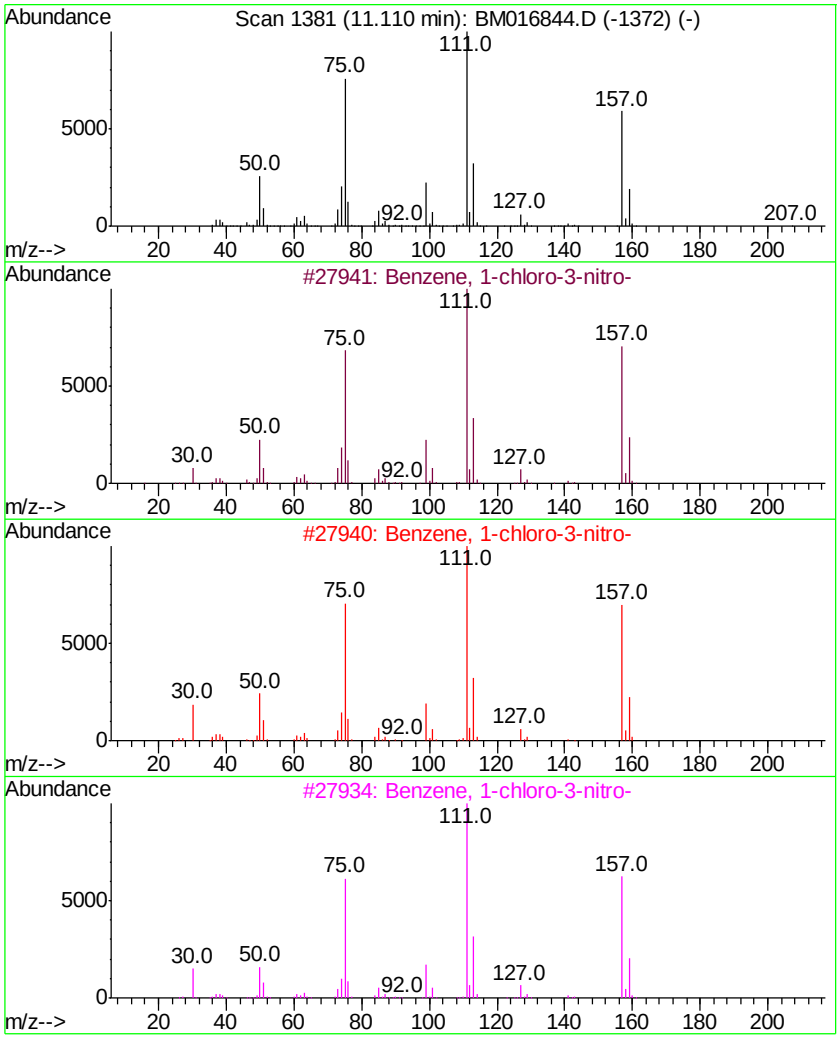
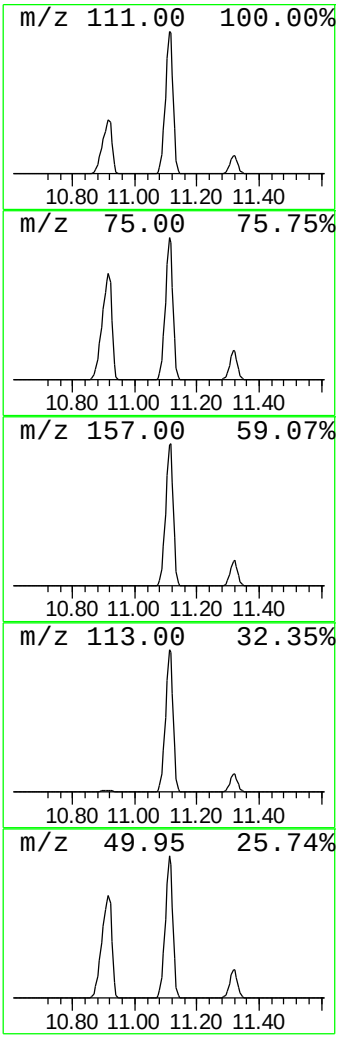
Peak Number 9 Benzene, 1-chloro-3-nitro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.11	152.80 ng/ul	17729900	Naphthalene-d8	10.53

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	98
2	Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	97
3	Benzene, 1-chloro-3-nitro-	157	C6H4ClNO2	000121-73-3	96
4	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	94
5	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	91





Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092118\
Data File : BM016844.D
Acq On : 21 Sep 2018 22:18
Operator : SJ/JU
Sample : J5013-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08R4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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

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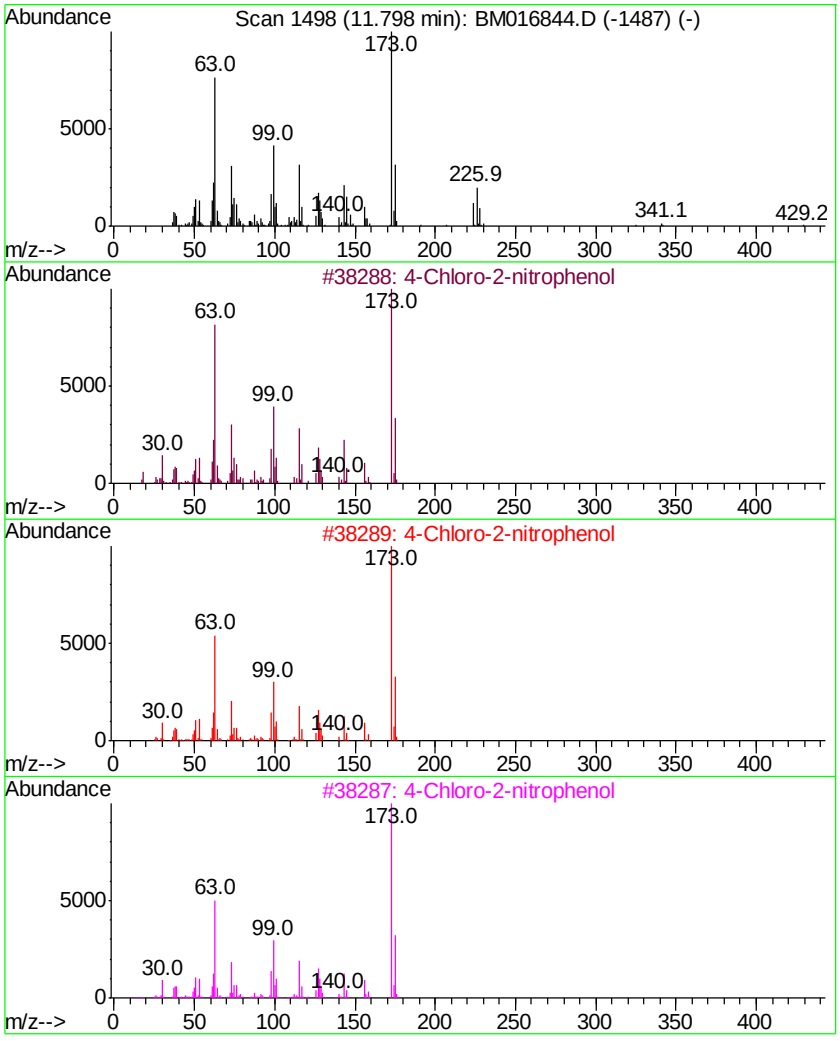
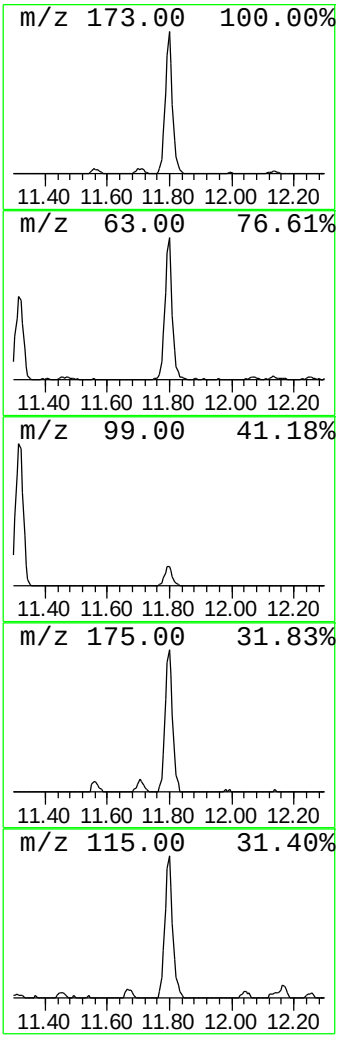
Peak Number 11 4-Chloro-2-nitrophenol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.80	7.49 ng/ul	869438	Naphthalene-d8	10.53

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	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, 5-chloro-2-nitro-	173	C6H4ClNO3	000611-07-4	87
5	3-Chloro-4-nitrophenol	173	C6H4ClNO3	000491-11-2	68





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Acq On : 21 Sep 2018 22:18
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Sample : J5013-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08R4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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

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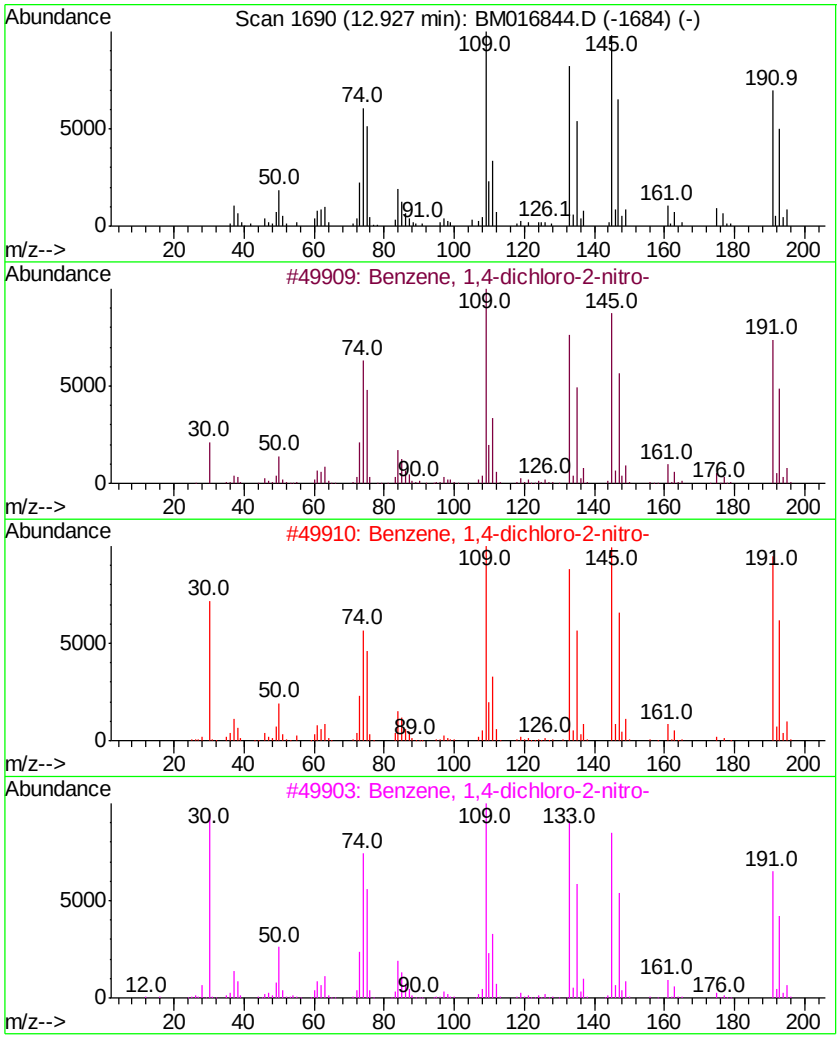
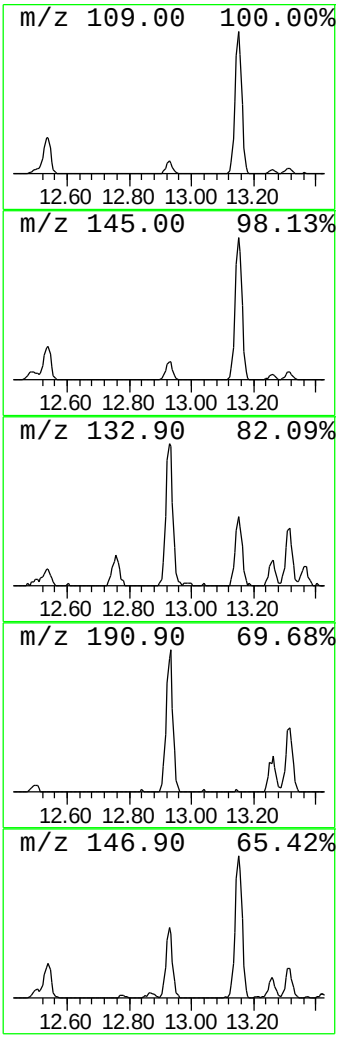
Peak Number 13 Benzene, 1,4-dichloro-2-nitro- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.93	2.69 ng/ul	401429	Acenaphthene-d10	14.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-dichloro-2-nitro-	191	C6H3Cl2NO2	000089-61-2	99
2	Benzene, 1,4-dichloro-2-nitro-	191	C6H3Cl2NO2	000089-61-2	99
3	Benzene, 1,4-dichloro-2-nitro-	191	C6H3Cl2NO2	000089-61-2	98
4	Benzene, 1,2-dichloro-3-nitro-	191	C6H3Cl2NO2	003209-22-1	93
5	Benzene, 2,4-dichloro-1-nitro-	191	C6H3Cl2NO2	000611-06-3	93





Library Search Compound Report

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Quant Title : SVOA CALIBRATION

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

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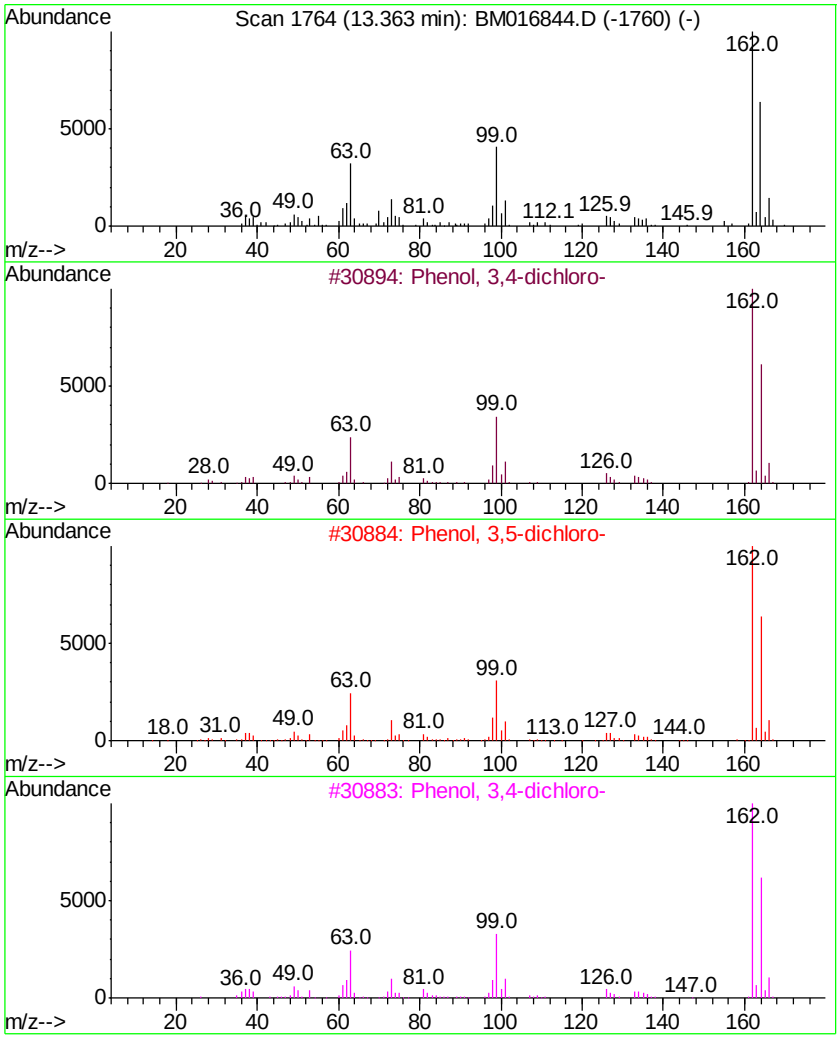
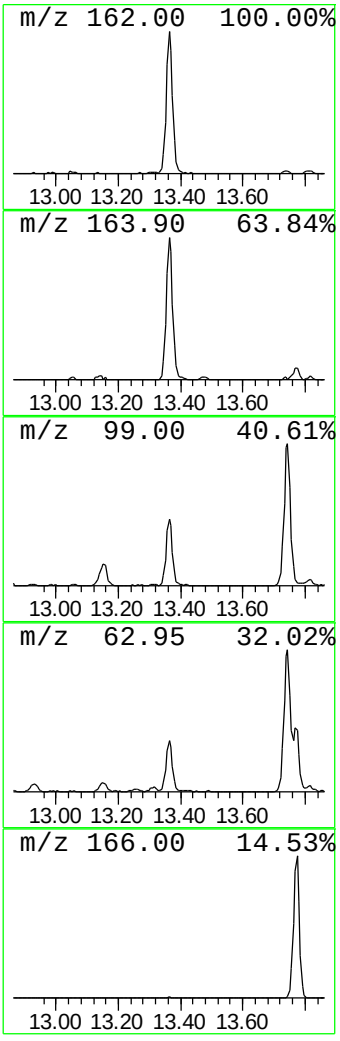
Peak Number 14 Phenol, 3,4-dichloro- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.36	2.35 ng/ul	351122	Acenaphthene-d10	14.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	98
2	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	98
3	Phenol, 3,4-dichloro-	162	C6H4Cl2O	000095-77-2	98
4	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	96
5	Phenol, 3,5-dichloro-	162	C6H4Cl2O	000591-35-5	96





Library Search Compound Report

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Data File : BM016844.D
Acq On : 21 Sep 2018 22:18
Operator : SJ/JU
Sample : J5013-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08R4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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

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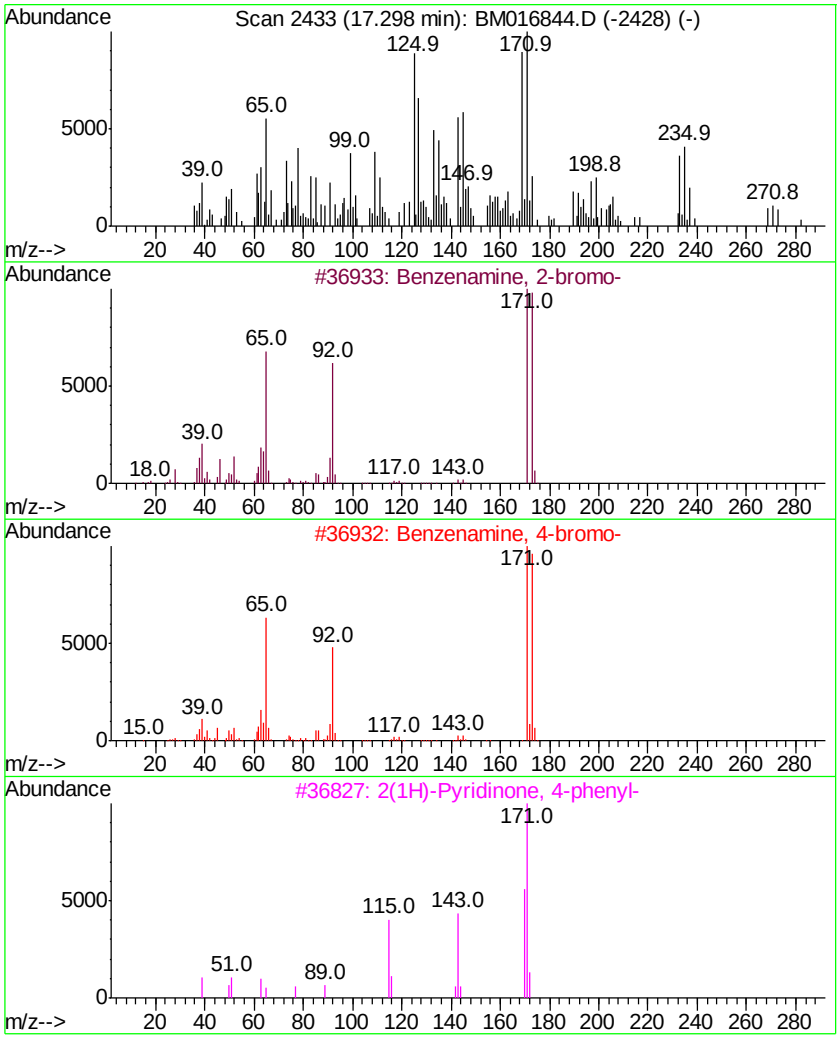
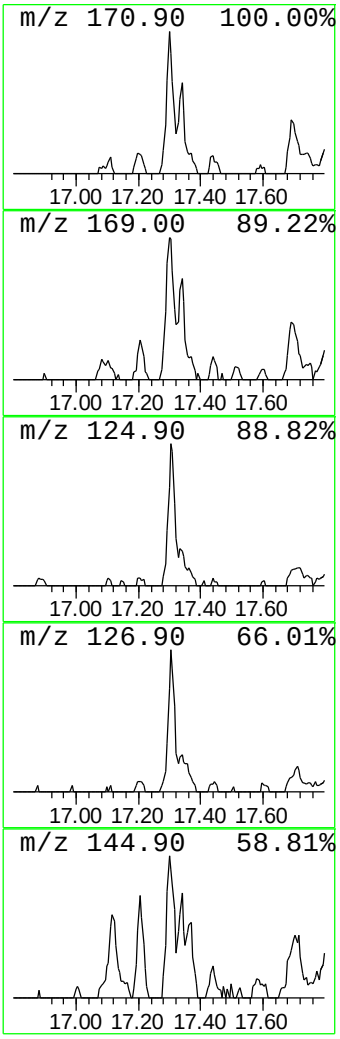
Peak Number 15 unknown-01 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.30	2.12 ng/ul	352958	Phenanthrene-d10	17.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzenamine, 2-bromo-	171	C6H6BrN	000615-36-1	27
2	Benzenamine, 4-bromo-	171	C6H6BrN	000106-40-1	16
3	2(1H)-Pyridinone, 4-phenyl-	171	C11H9NO	019006-81-6	11
4	p-Bromo-benzyl tetrahydropyran-2...	270	C12H15BrO2	017100-68-4	10
5	Formamide, N-1-naphthalenyl-	171	C11H9NO	006330-51-4	10





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Data File : BM016844.D
Acq On : 21 Sep 2018 22:18
Operator : SJ/JU
Sample : J5013-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08R4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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

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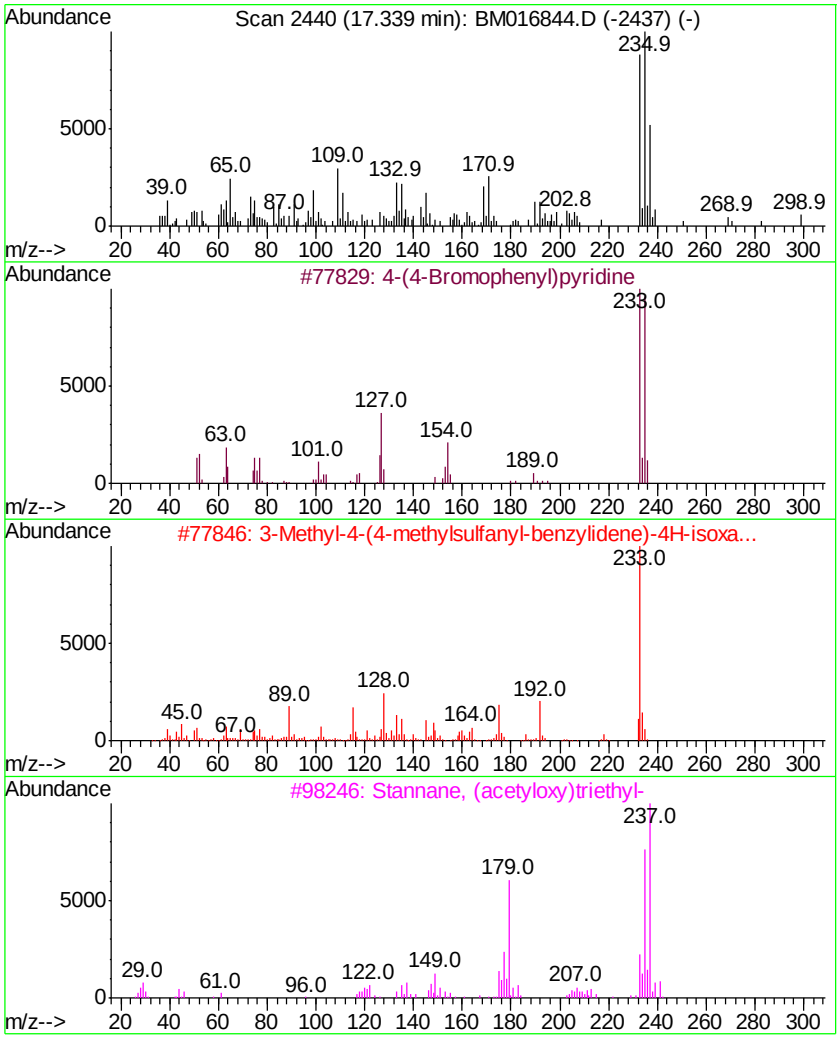
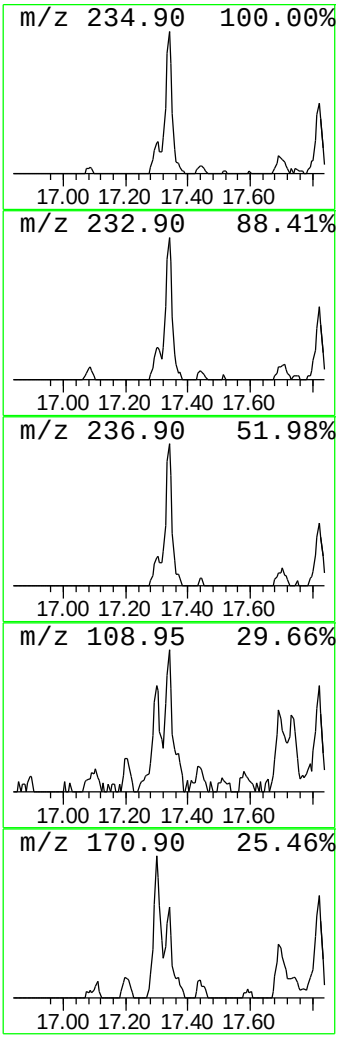
Peak Number 16 unknown-02 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.34	2.17 ng/ul	360838	Phenanthrene-d10	17.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-(4-Bromophenyl)pyridine	233	C11H8BrN	039795-60-3	35
2	3-Methyl-4-(4-methylsulfanyl-ben...	233	C12H11NO2S	146795-09-7	22
3	Stannane, (acetyloxy)triethyl-	266	C8H18O2Sn	001907-13-7	22
4	3-Chloro-7-ethyl-1-(3-hydroxy-pr...	294	C14H19ClN4O	1000274-38-9	22
5	4-Isothiazolecarboxylic acid, 5-...	233	C12H11NO2S	021905-49-7	18





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Data File : BM016844.D
Acq On : 21 Sep 2018 22:18
Operator : SJ/JU
Sample : J5013-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08R4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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

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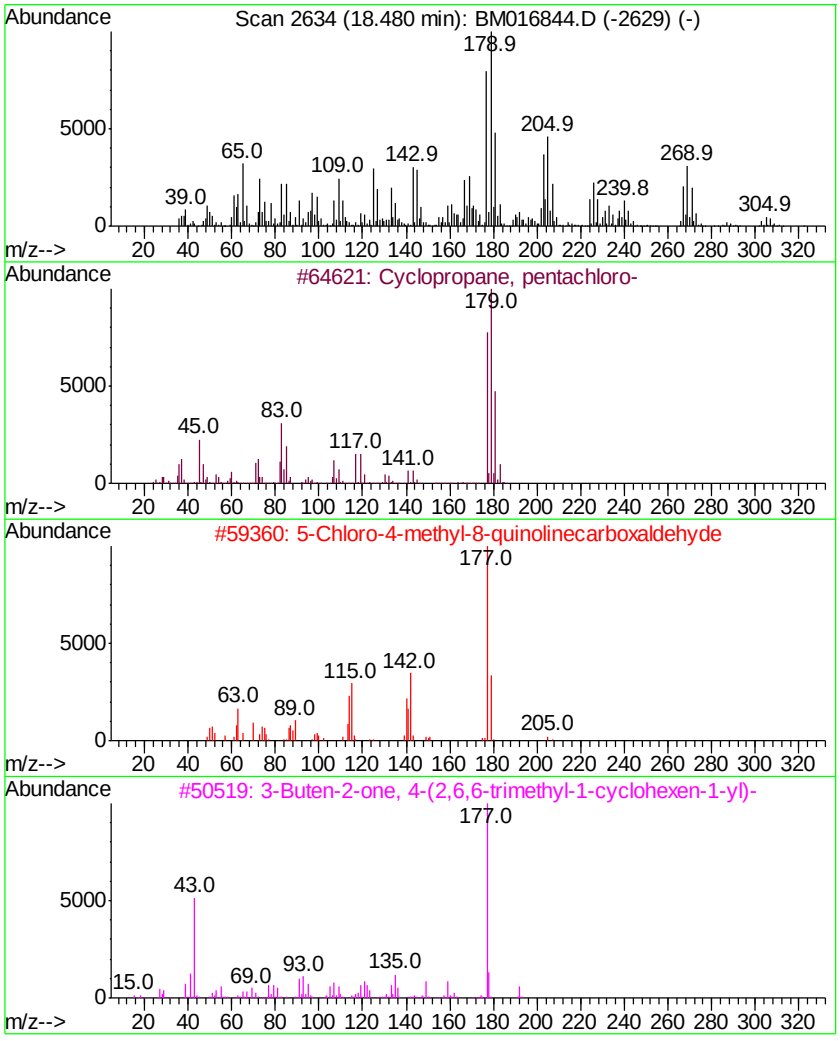
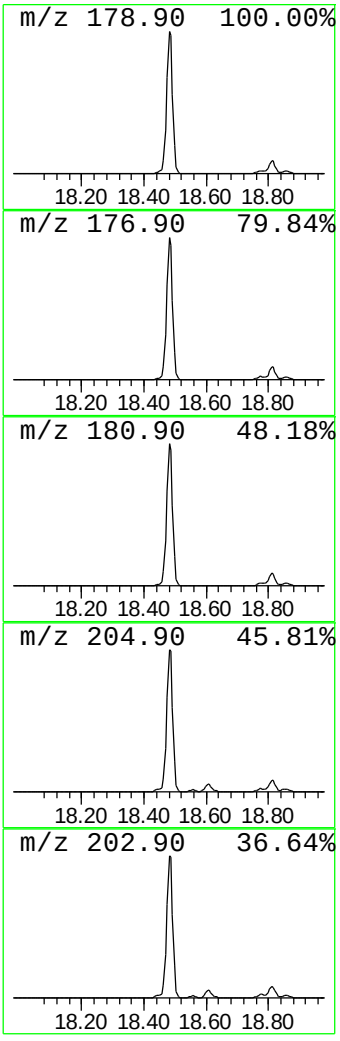
Peak Number 17 unknown-03 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.48	103.82 ng/ul	17294100	Phenanthrene-d10	17.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, pentachloro-	212	C3HCl5	006262-51-7	43
2	5-Chloro-4-methyl-8-quinolinecar...	205	C11H8ClNO	028662-47-7	27
3	3-Buten-2-one, 4-(2,6,6-trimethy...	192	C13H20O	014901-07-6	25
4	Octamethyl-1,4-cyclohexadiene	192	C14H24	172684-93-4	25
5	2,4,6-Trifluoronitrobenzene	177	C6H2F3NO2	000315-14-0	18





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092118\
Data File : BM016844.D
Acq On : 21 Sep 2018 22:18
Operator : SJ/JU
Sample : J5013-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08R4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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

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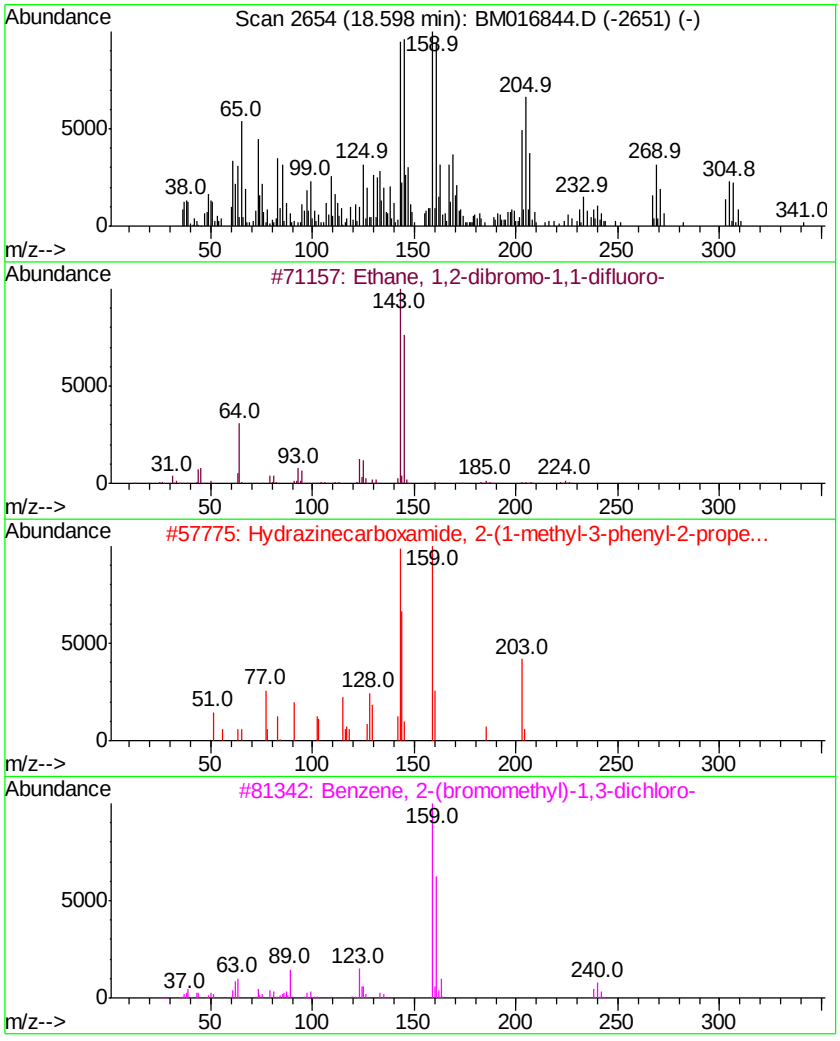
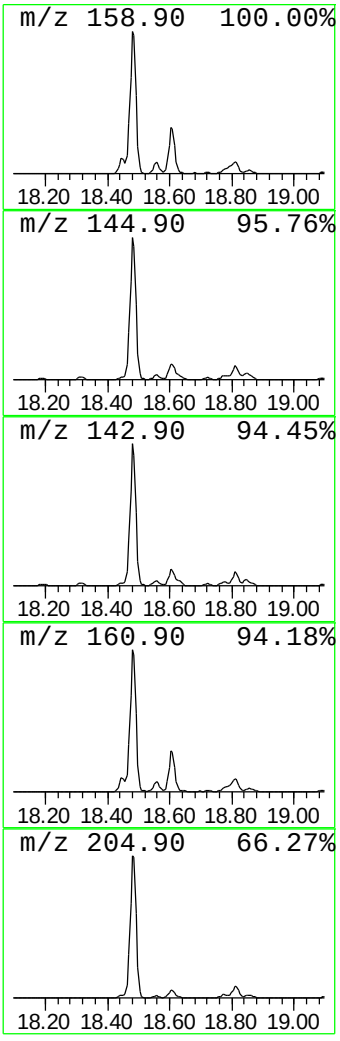
Peak Number 18 unknown-04 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.60	6.02 ng/ul	1002830	Phenanthrene-d10	17.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethane, 1,2-dibromo-1,1-difluoro-	222	C2H2Br2F2	000075-82-1	11
2	Hydrazinecarboxamide, 2-(1-methy...	203	C11H13N3O	005468-31-5	10
3	Benzene, 2-(bromomethyl)-1,3-dic...	238	C7H5BrCl2	020443-98-5	10
4	Benzo[b]furan-3-carboxamide	161	C9H7NO2	1000262-22-7	10
5	Benzene, 2-(bromomethyl)-1,3-dic...	238	C7H5BrCl2	020443-98-5	10





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Data File : BM016844.D
Acq On : 21 Sep 2018 22:18
Operator : SJ/JU
Sample : J5013-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08R4

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM091018MA.M
Quant Title : SVOA CALIBRATION

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

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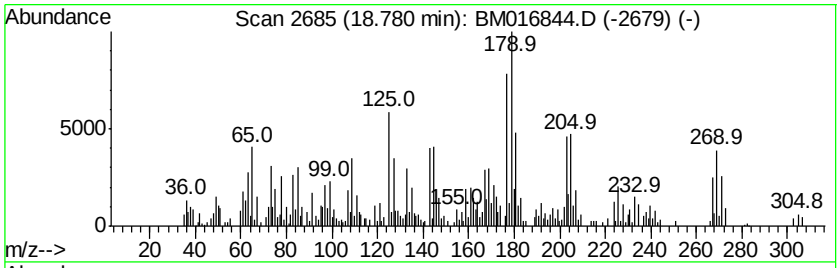
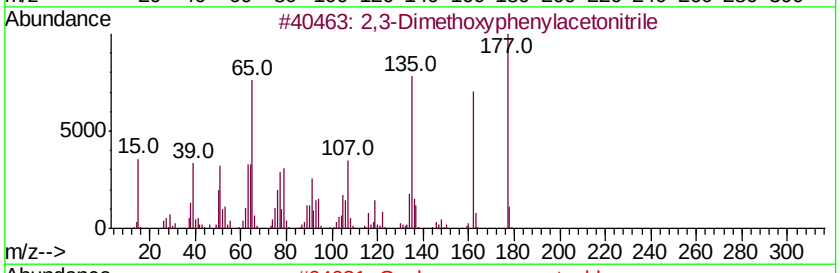
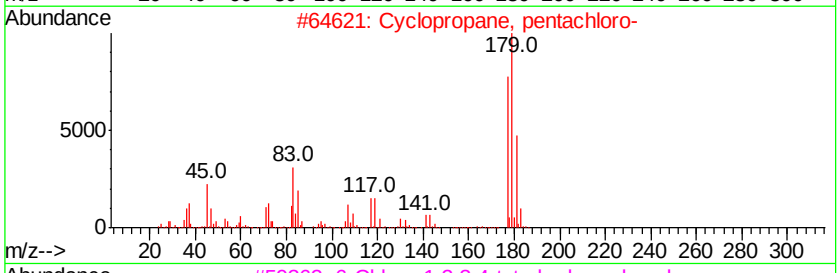
Peak Number 19 unknown-05 Concentration Rank 15

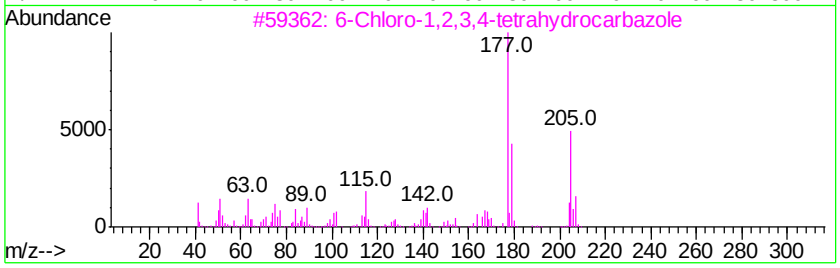
R.T.	EstConc	Area	Relative to ISTD	R.T.
18.78	3.15 ng/ul	525562	Phenanthrene-d10	17.10

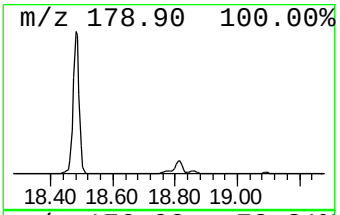
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3-Dimethoxyphenylacetonitrile	177	C10H11NO2	004468-57-9	27
2	Cyclopropane, pentachloro-	212	C3HCl5	006262-51-7	12
3	6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	11
4	Trimethylphenylgermanium	196	C9H14Ge	001626-00-2	11
5	Isothiazole, 5-bromo-3-methyl-	177	C4H4BrNS	020493-60-1	9

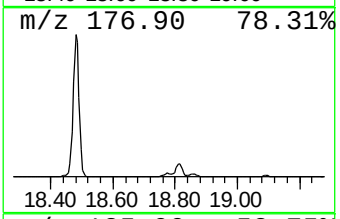


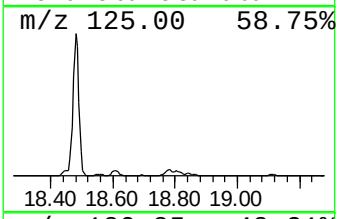


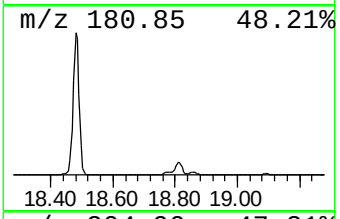


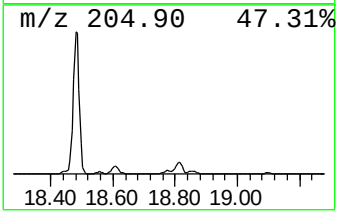












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Misc :
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Instrument : BNA_M
ClientSampleId : C08R4

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Quant Title : SVOA CALIBRATION

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

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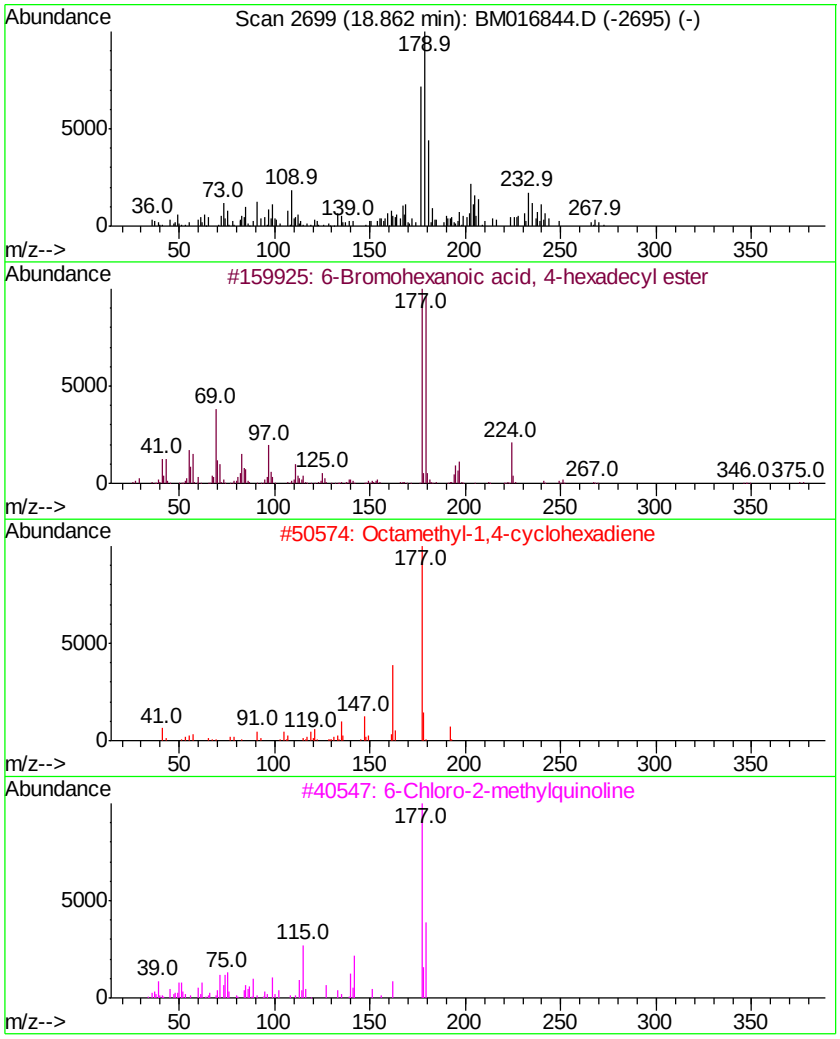
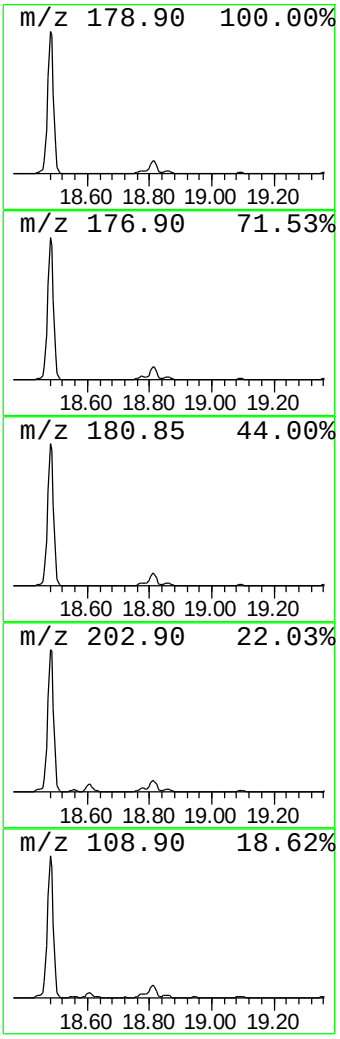
Peak Number 21 unknown-06 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.86	2.33 ng/ul	387557	Phenanthrene-d10	17.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Bromohexanoic acid, 4-hexadecv...	418	C22H43BrO2	1000282-90-6	25
2	Octamethyl-1,4-cyclohexadiene	192	C14H24	172684-93-4	18
3	6-Chloro-2-methylquinoline	177	C10H8ClN	000092-46-6	16
4	4(1H)-Pteridinone, 2-amino-7-met...	177	C7H7N5O	013040-58-9	14
5	Benzothiazole-6-carboxylic acid	179	C8H5NO2S	003622-35-3	14





Tentatively Identified Compound (LSC) summary

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM092118\
Data File : BM016844.D
Acq On : 21 Sep 2018 22:18
Operator : SJ/JU
Sample : J5013-01
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08R4

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Quant Title : SVOA CALIBRATION

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, 1-bromo-...	9.22	20.8	ng/ul	2408440	2	10.53	2320700	20.0
Benzene, 1,2-dich...	9.44	2.3	ng/ul	261751	2	10.53	2320700	20.0
Benzene, 1-chloro...	11.11	152.8	ng/ul	17729900	2	10.53	2320700	20.0
4-Chloro-2-nitrop...	11.80	7.5	ng/ul	869438	2	10.53	2320700	20.0
Benzene, 1,4-dich...	12.93	2.7	ng/ul	401429	3	14.36	2983260	20.0
Phenol, 3,4-dichl...	13.36	2.4	ng/ul	351122	3	14.36	2983260	20.0
unknown-01	17.30	2.1	ng/ul	352958	4	17.10	3331710	20.0
unknown-02	17.34	2.2	ng/ul	360838	4	17.10	3331710	20.0
unknown-03	18.48	103.8	ng/ul	17294100	4	17.10	3331710	20.0
unknown-04	18.60	6.0	ng/ul	1002830	4	17.10	3331710	20.0
unknown-05	18.78	3.1	ng/ul	525562	4	17.10	3331710	20.0
unknown-06	18.86	2.3	ng/ul	387557	4	17.10	3331710	20.0