LSC Area Percent Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092118\
Data File : BM016839.D
Acq On : 21 Sep 2018 19:17
Operator : SJ/JU
Sample : J5012-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08K5

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	2565033	2956639	1.52%	0.274%
2	3.987	165	170	178	rVB	124126	209353	0.11%	0.019%
3	5.093	348	358	361	rBV4	50802062	127554214	65.43%	11.842%
4	6.422	577	584	594	rVB	789364	1302311	0.67%	0.121%
5	6.687	611	629	632	rBV	290025	585648	0.30%	0.054%
6	6.899	658	665	677	rVB	271951	522457	0.27%	0.049%
7	7.099	687	699	708	rBV	505773	1606196	0.82%	0.149%
8	7.269	721	728	738	rVB	742723	1724062	0.88%	0.160%
9	7.469	754	762	774	rVB3	81670	312436	0.16%	0.029%
10	7.640	779	791	806	rBV2	28770962	144770241	74.26%	13.440%
11	7.798	806	818	820	rBV3	53097873	159400485	81.76%	14.798%
12	8.140	861	876	879	rBV4	54102727	194771416	99.90%	18.082%
13	8.434	919	926	935	rVB	887039	1376798	0.71%	0.128%
14	8.893	996	1004	1006	rBV	1114569	1990298	1.02%	0.185%
15	8.945	1006	1013	1019	rVB	13986846	25593063	13.13%	2.376%
16	9.216	1052	1059	1067	rBV	1192992	1941026	1.00%	0.180%
17	9.440	1091	1097	1101	rBV	139057	231569	0.12%	0.021%
18	9.487	1101	1105	1107	rVV	125934	203044	0.10%	0.019%
19	9.528	1107	1112	1118	rVV	601102	1019406	0.52%	0.095%
20	9.598	1118	1124	1139	rVB2	1014011	2299714	1.18%	0.213%
21	10.128	1207	1214	1225	rBV	1490475	2559491	1.31%	0.238%
22	10.416	1247	1263	1268	rBV5	57832426	194958600	100.00%	18.099%
23	10.522	1275	1281	1287	rBV	1603797	2443540	1.25%	0.227%
24	10.910	1335	1347	1353	rBV	41624862	83772690	42.97%	7.777%
25	11.116	1372	1382	1393	rBV	12597088	21642819	11.10%	2.009%
26	11.322	1408	1417	1425	rVB	2710672	4714723	2.42%	0.438%
27	11.798	1488	1498	1508	rBV2	426458	776940	0.40%	0.072%
28	12.486	1607	1615	1619	rBV	504334	1060763	0.54%	0.098%
29	12.539	1619	1624	1631	rVB	2578905	4129010	2.12%	0.383%
30	12.763	1655	1662	1674	rVB2	132574	332643	0.17%	0.031%
31	12.928	1684	1690	1696	rBV	325890	486021	0.25%	0.045%
32	13.151	1721	1728	1737	rVB	7918394	11731050	6.02%	1.089%
33	13.316	1751	1756	1761	rVV2	137977	225857	0.12%	0.021%
34	13.745	1822	1829	1830	rBV	804240	1103508	0.57%	0.102%

LSC Area Percent Report

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Data File : BM016839.D
Acq On : 21 Sep 2018 19:17
Operator : SJ/JU
Sample : J5012-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08K5

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	13.775	1830	1834	1839	rVV	2805447	3915610	2.01%	0.364%
36	13.992	1865	1871	1875	rBV3	164026	275701	0.14%	0.026%
37	14.051	1875	1881	1890	rVV	3398898	4964649	2.55%	0.461%
38	14.357	1928	1933	1947	rVB2	1998288	3190729	1.64%	0.296%
39	14.551	1962	1966	1976	rVB2	308745	522064	0.27%	0.048%
40	14.845	2010	2016	2021	rBV	125333	200013	0.10%	0.019%
41	15.357	2095	2103	2115	rVB	4156732	6184388	3.17%	0.574%
42	15.469	2115	2122	2135	rBV	1155955	1608345	0.82%	0.149%
43	17.104	2393	2400	2408	rBV	2570456	3596682	1.84%	0.334%
44	17.204	2410	2417	2428	rVV	4618924	6363156	3.26%	0.591%
45	18.486	2629	2635	2643	rVB	12564175	17622271	9.04%	1.636%
46	18.557	2643	2647	2651	rVV3	163604	227430	0.12%	0.021%
47	18.604	2651	2655	2665	rVB	483670	732702	0.38%	0.068%
48	18.780	2679	2685	2687	rBV3	323904	562469	0.29%	0.052%
49	18.815	2687	2691	2695	rVV	881541	1317484	0.68%	0.122%
50	18.857	2695	2698	2703	rVB3	244636	379983	0.19%	0.035%
51	19.098	2733	2739	2743	rBV4	108009	216418	0.11%	0.020%
52	19.362	2780	2784	2794	rVB2	265191	419318	0.22%	0.039%
53	19.504	2801	2808	2816	rBV	5397818	7569361	3.88%	0.703%
54	21.292	3107	3112	3119	rBV	3529365	4367887	2.24%	0.405%
55	23.386	3461	3468	3475	rBV	4421239	8146392	4.18%	0.756%
56	23.527	3486	3492	3504	rVB2	2361771	4484439	2.30%	0.416%

Sum of corrected areas: 1077175522

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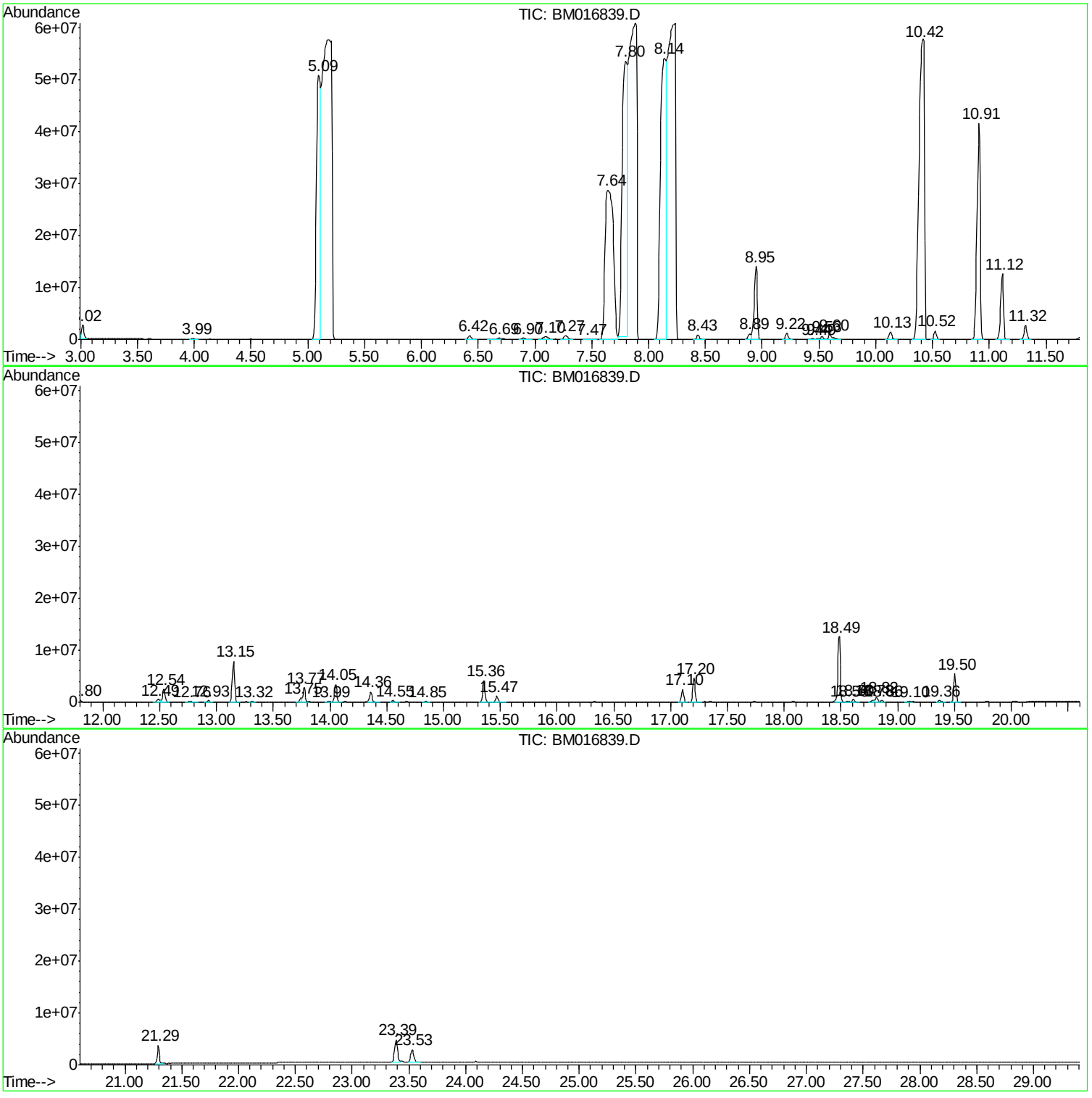
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C08K5

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

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Data File : BM016839.D
Acq On : 21 Sep 2018 19:17
Operator : SJ/JU
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Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08K5

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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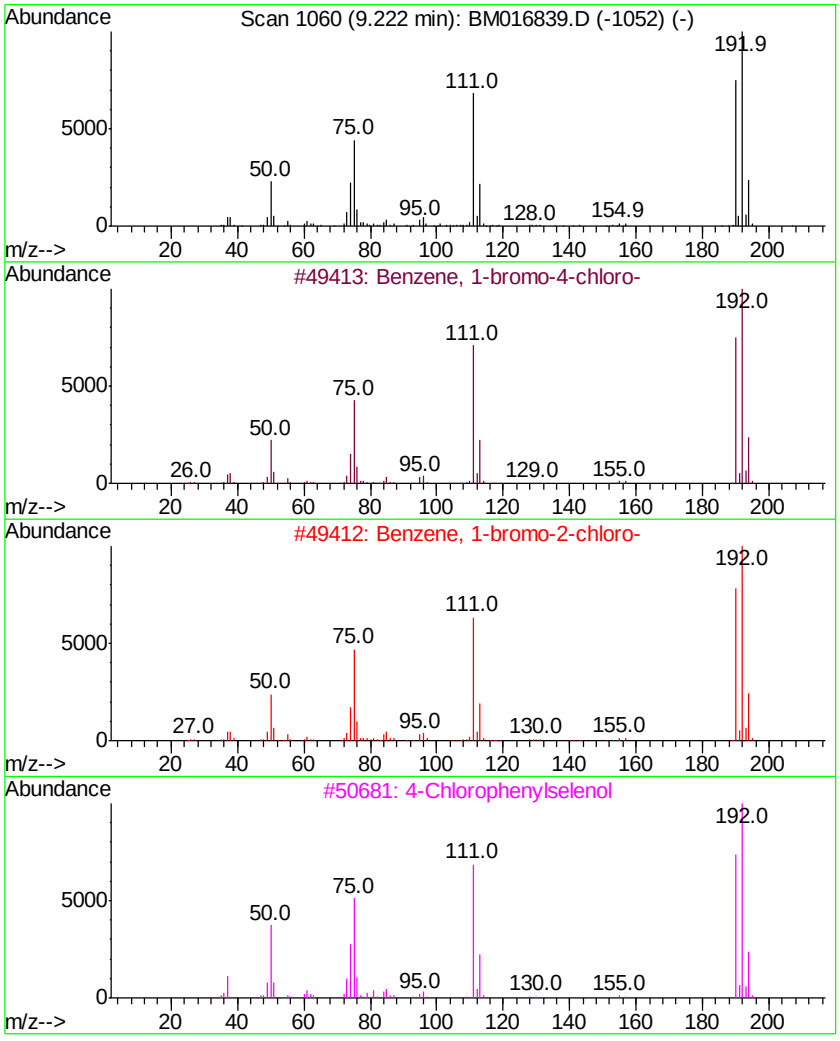
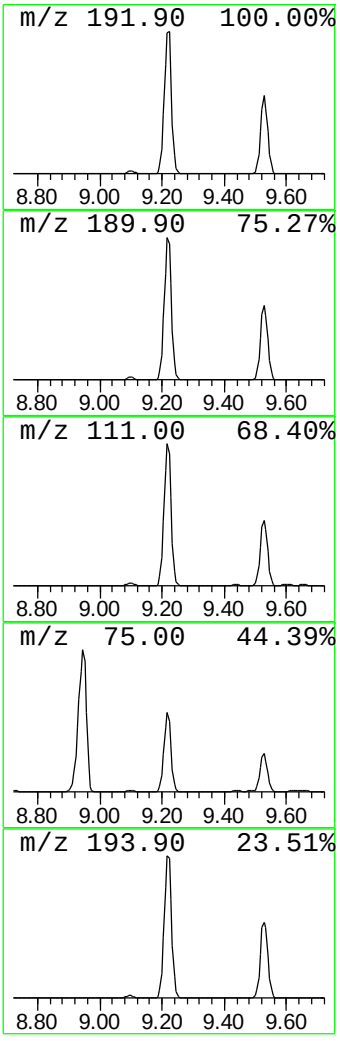
Peak Number 4 Benzene, 1-bromo-4-chloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.22	15.89 ng/ul	1941030	Naphthalene-d8	10.52

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	98
2	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
3	4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	97
4	Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	97
5	Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	96





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092118\
Data File : BM016839.D
Acq On : 21 Sep 2018 19:17
Operator : SJ/JU
Sample : J5012-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08K5

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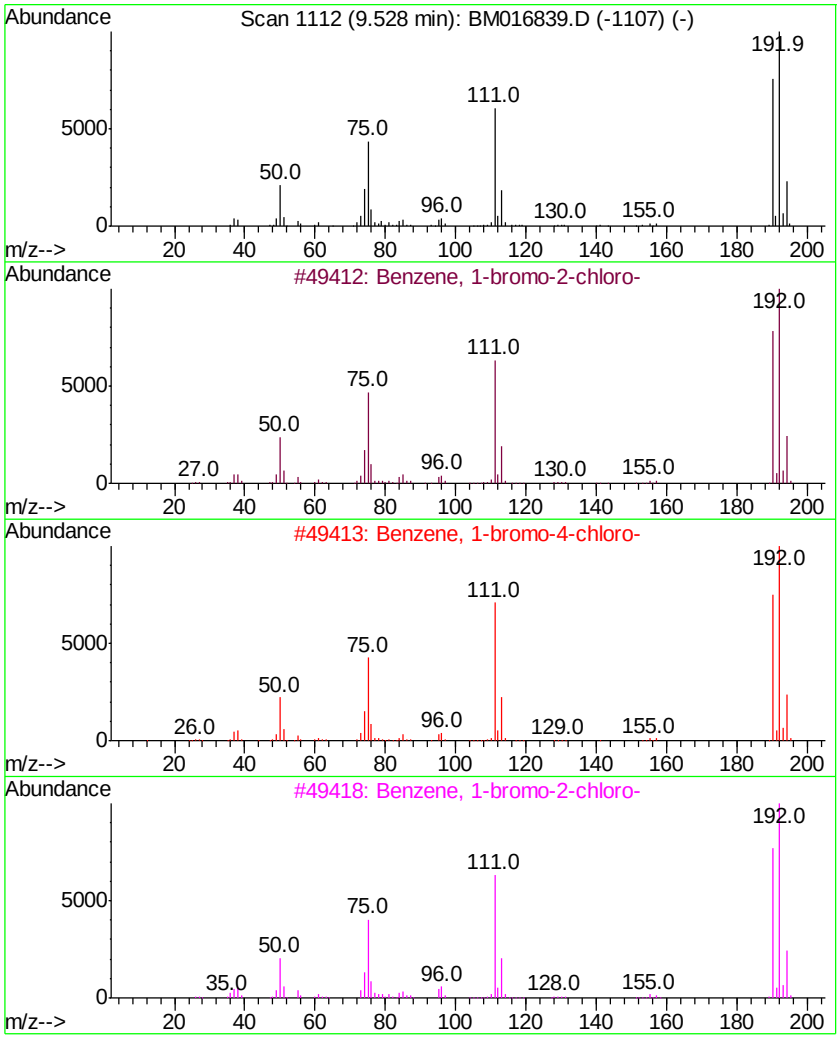
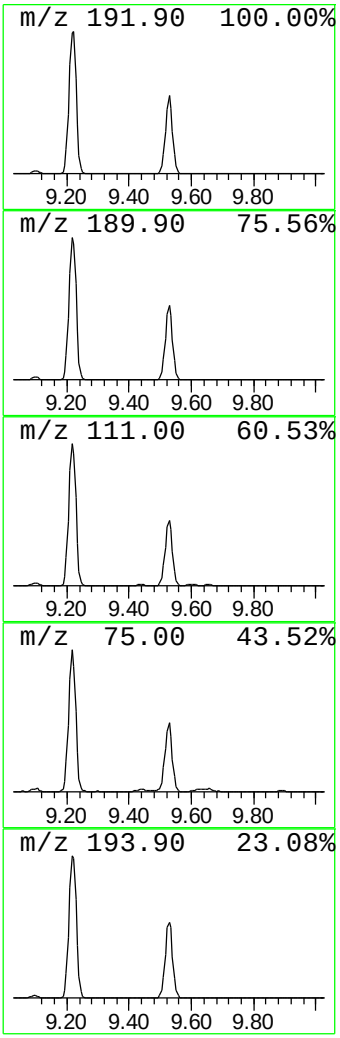
Peak Number 5 Benzene, 1-bromo-2-chloro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.53	8.34 ng/ul	1019410	Naphthalene-d8	10.52

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	Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	96





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092118\
Data File : BM016839.D
Acq On : 21 Sep 2018 19:17
Operator : SJ/JU
Sample : J5012-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08K5

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

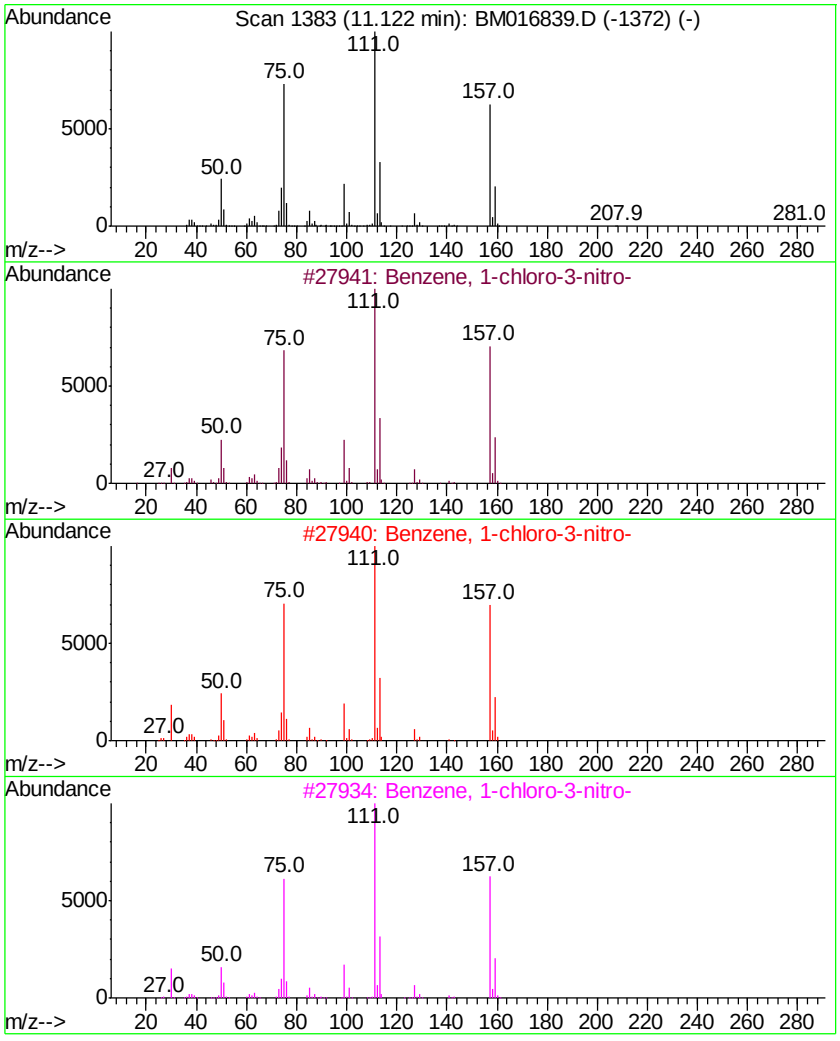
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

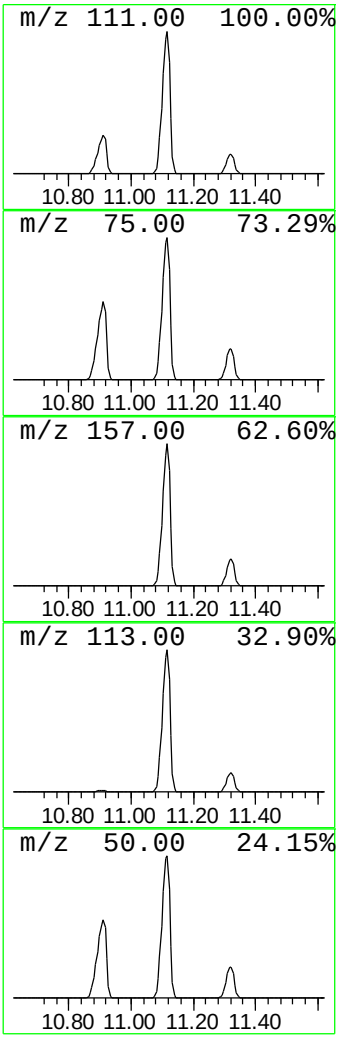
****

Peak Number 8 Benzene, 1-chloro-3-nitro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.12	177.14 ng/ul	21642800	Naphthalene-d8	10.52

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	91
5	Benzene, 1-chloro-4-nitro-	157	C6H4ClNO2	000100-00-5	90





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092118\
Data File : BM016839.D
Acq On : 21 Sep 2018 19:17
Operator : SJ/JU
Sample : J5012-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08K5

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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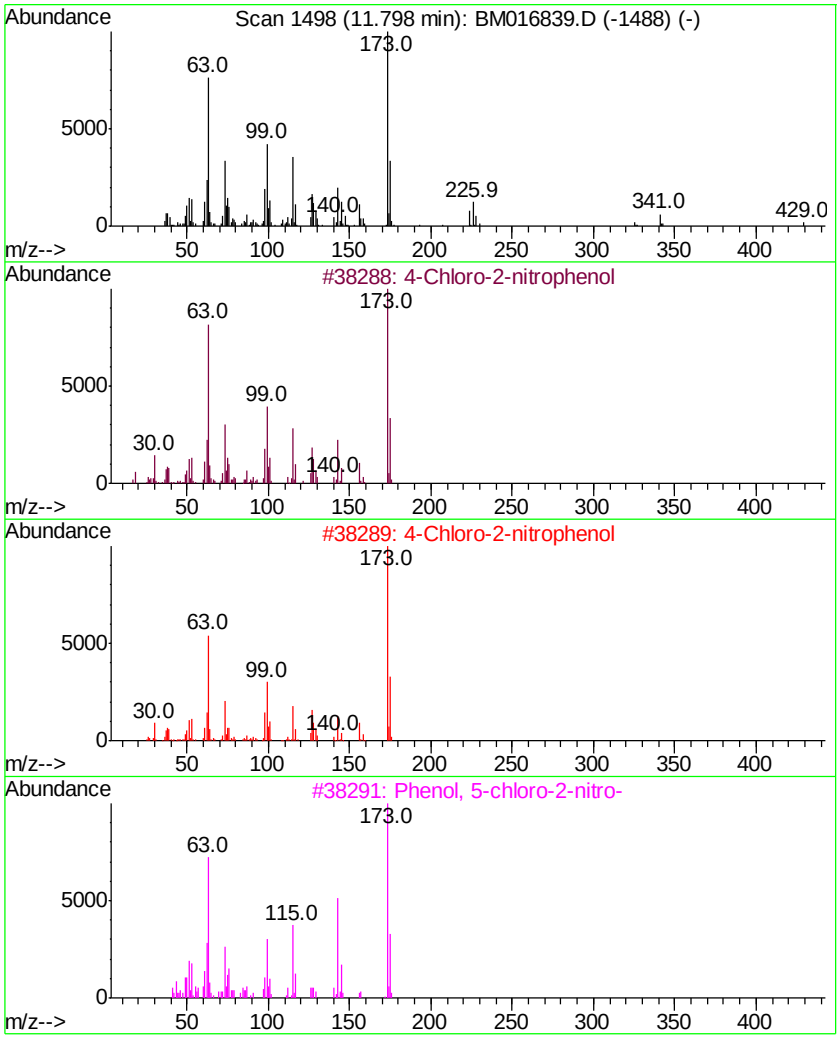
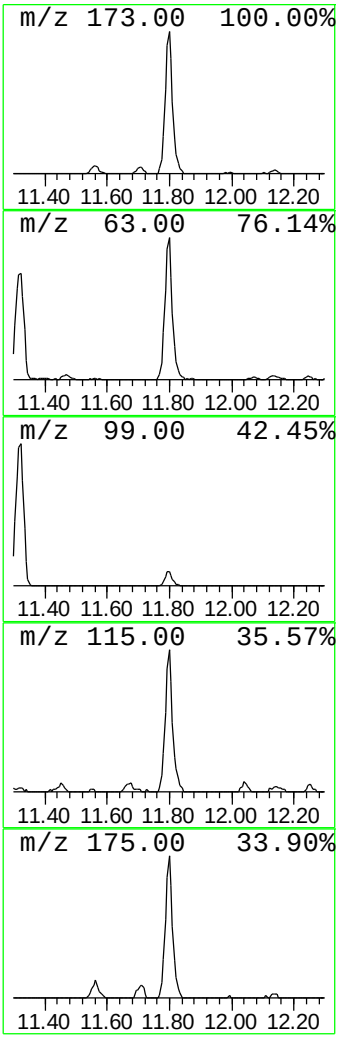
Peak Number 10 4-Chloro-2-nitrophenol Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.80	6.36 ng/ul	776940	Naphthalene-d8	10.52

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	93
3	Phenol, 5-chloro-2-nitro-	173	C6H4ClNO3	000611-07-4	87
4	4-Chloro-2-nitrophenol	173	C6H4ClNO3	000089-64-5	60
5	Propane, 1,1,2-trichloro-	146	C3H5Cl3	000598-77-6	50





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Instrument : BNA_M
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Quant Title : SVOA CALIBRATION

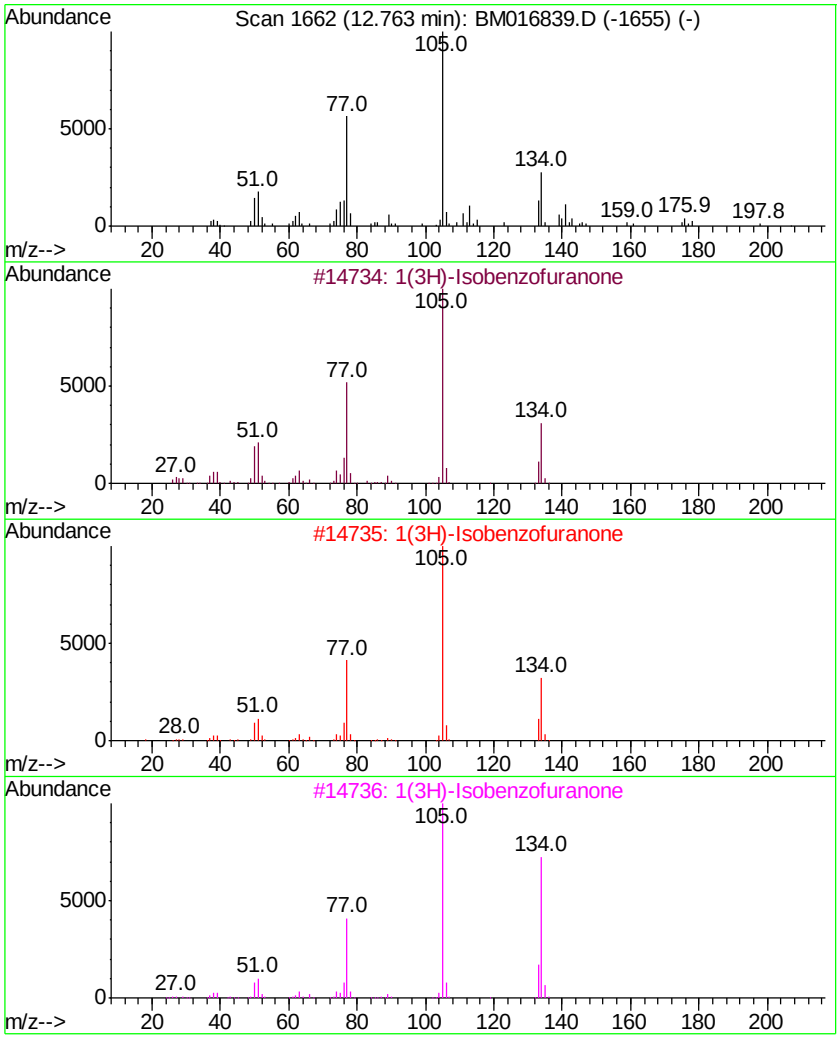
TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

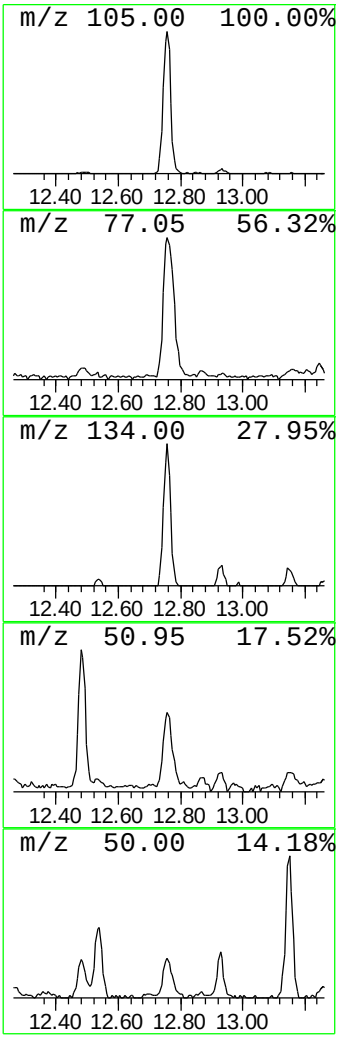
****

Peak Number 12 1(3H)-Isobenzofuranone Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.76	2.09 ng/ul	332643	Acenaphthene-d10	14.36

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1(3H)-Isobenzofuranone	134	C8H6O2	000087-41-2	93
2	1(3H)-Isobenzofuranone	134	C8H6O2	000087-41-2	76
3	1(3H)-Isobenzofuranone	134	C8H6O2	000087-41-2	76
4	1,2-Benzenedicarboxaldehyde	134	C8H6O2	000643-79-8	68
5	1,2-Benzenedicarboxaldehyde	134	C8H6O2	000643-79-8	50





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Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08K5

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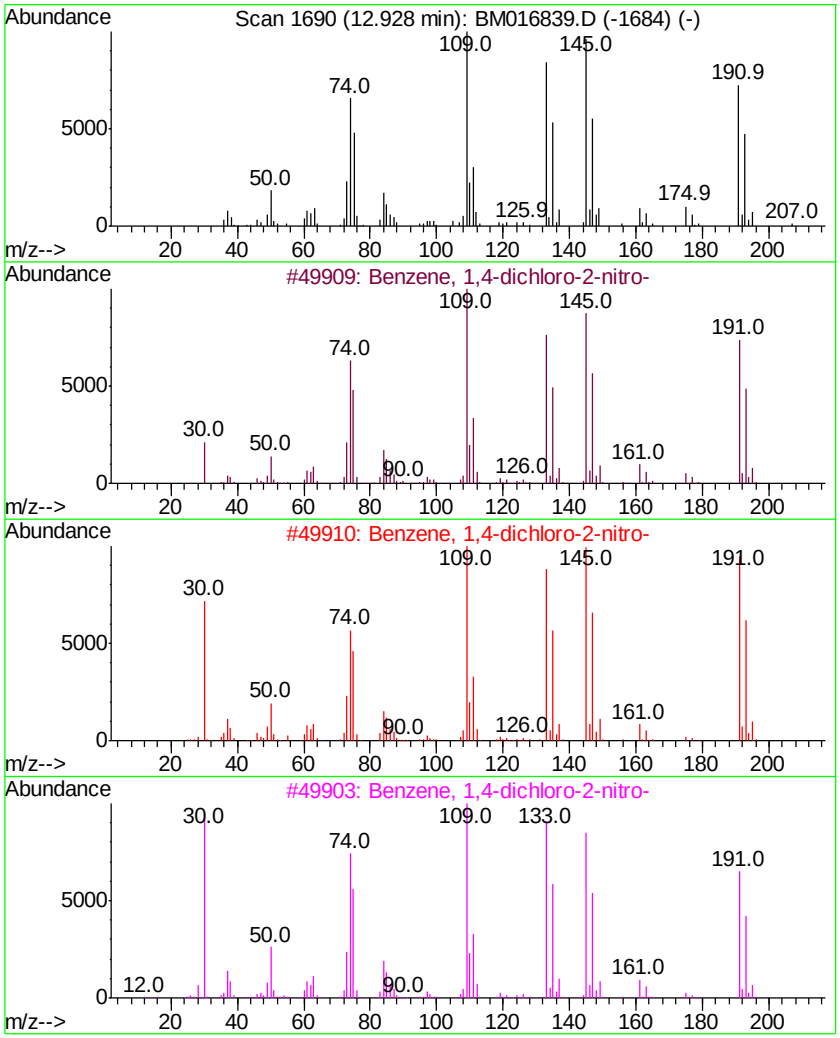
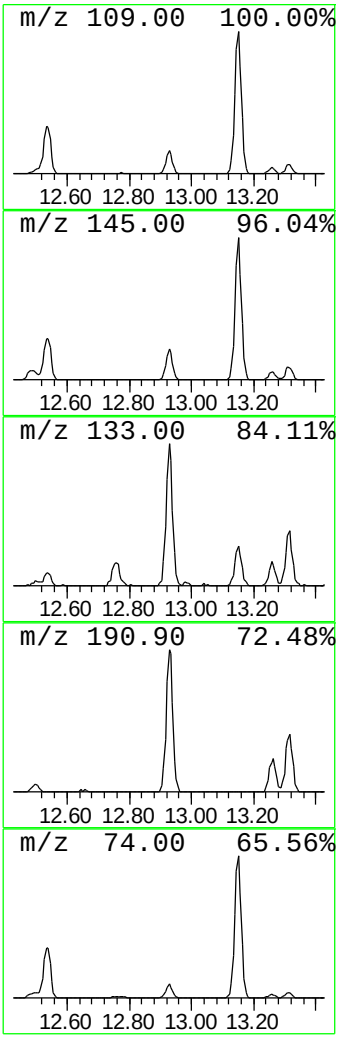
Peak Number 13 Benzene, 1,4-dichloro-2-nitro- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.93	3.05 ng/ul	486021	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	97
4	Benzene, 1,2-dichloro-3-nitro-	191	C6H3Cl2NO2	003209-22-1	94
5	Benzene, 2,4-dichloro-1-nitro-	191	C6H3Cl2NO2	000611-06-3	94





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Sample : J5012-03
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Instrument : BNA_M
ClientSampleId : C08K5

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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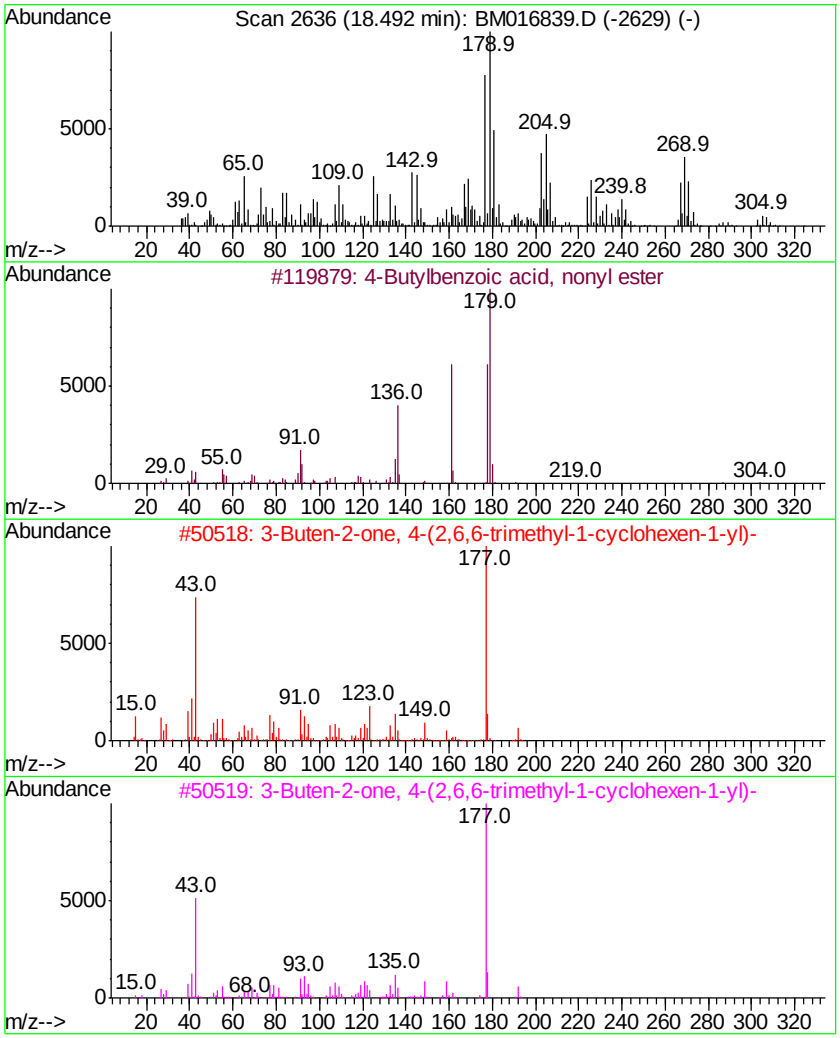
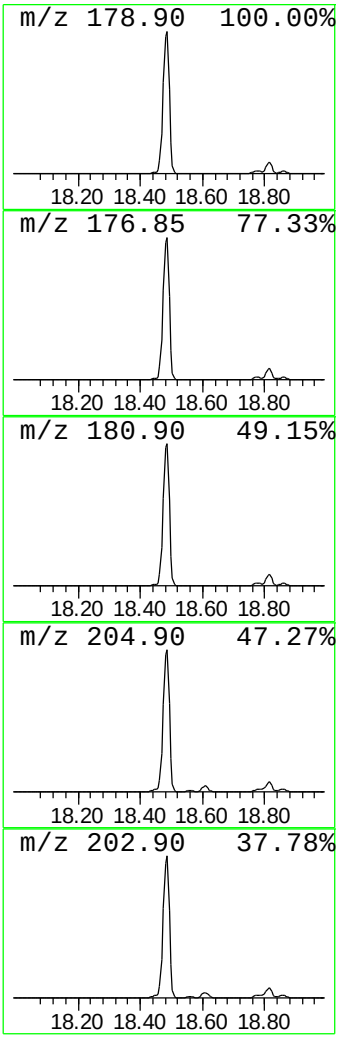
Peak Number 14 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.49	97.99 ng/ul	17622300	Phenanthrene-d10	17.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Butylbenzoic acid, nonyl ester	304	C20H32O2	1000292-20-6	25
2	3-Buten-2-one, 4-(2,6,6-trimethy...	192	C13H20O	014901-07-6	15
3	3-Buten-2-one, 4-(2,6,6-trimethy...	192	C13H20O	014901-07-6	15
4	5-Chloro-4-methyl-8-quinolinecar...	205	C11H8ClNO	028662-47-7	14
5	6-Bromohexanoic acid, 4-hexadecy...	418	C22H43BrO2	1000282-90-6	12





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Instrument : BNA_M
ClientSampleId : C08K5

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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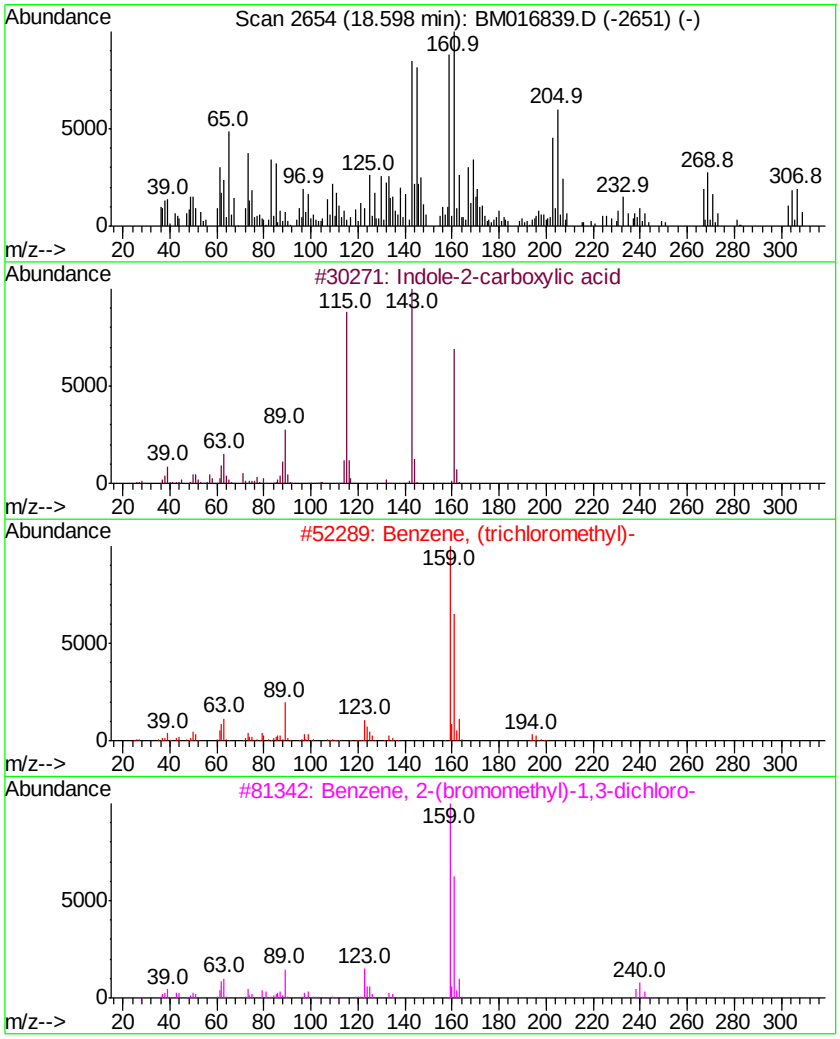
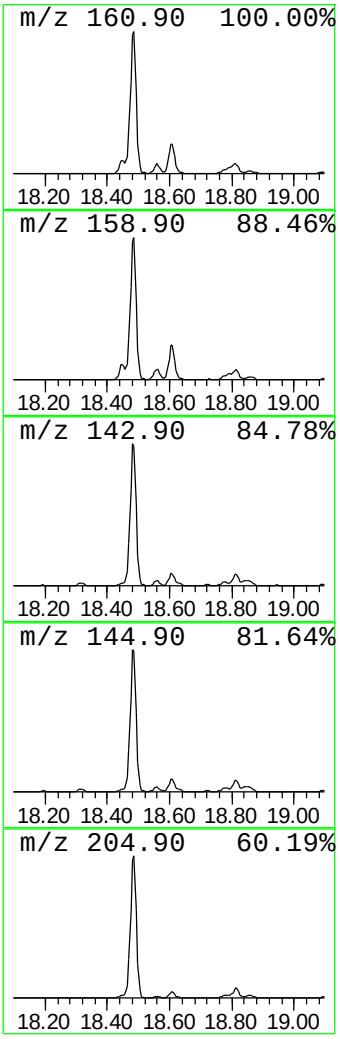
Peak Number 15 unknown-02 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.60	4.07 ng/ul	732702	Phenanthrene-d10	17.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indole-2-carboxylic acid	161	C9H7NO2	001477-50-5	18
2	Benzene, (trichloromethyl)-	194	C7H5Cl3	000098-07-7	18
3	Benzene, 2-(bromomethyl)-1,3-dic...	238	C7H5BrCl2	020443-98-5	14
4	Benzo[b]furan-3-carboxamide	161	C9H7NO2	1000262-22-7	11
5	Acetamide, 2-(2-ethyl-1-benzimid...	203	C11H13N3O	054980-94-8	11





Library Search Compound Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092118\
Data File : BM016839.D
Acq On : 21 Sep 2018 19:17
Operator : SJ/JU
Sample : J5012-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08K5

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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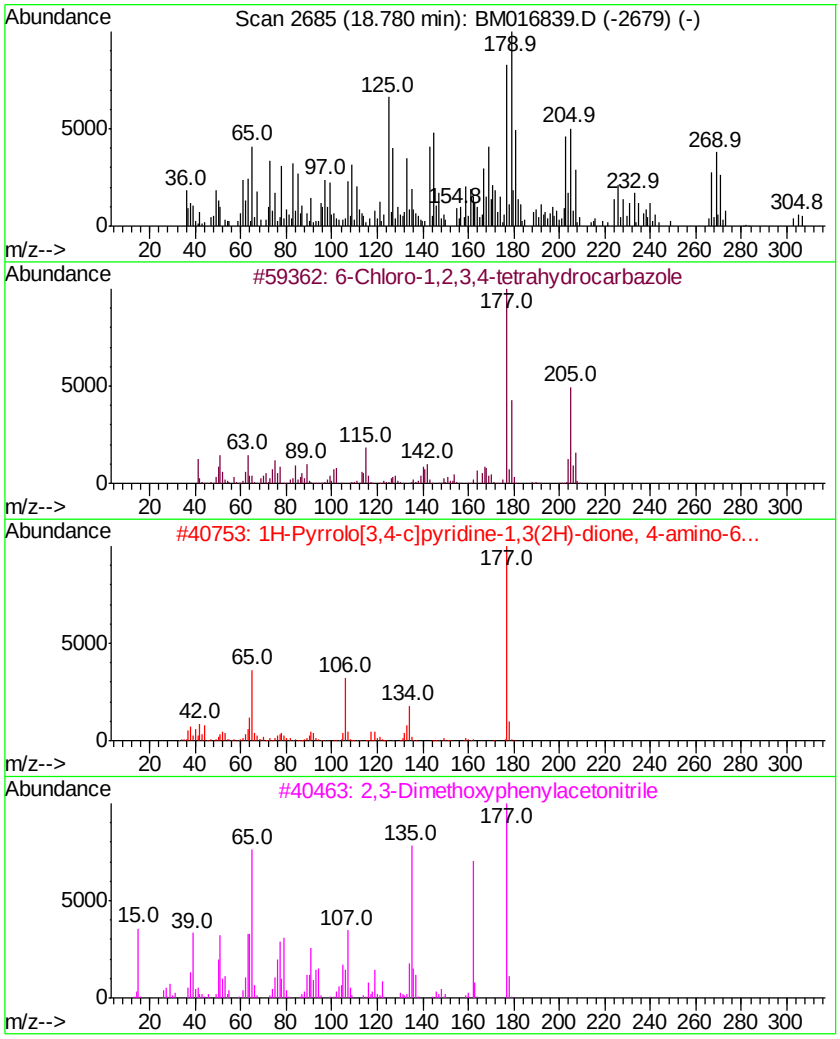
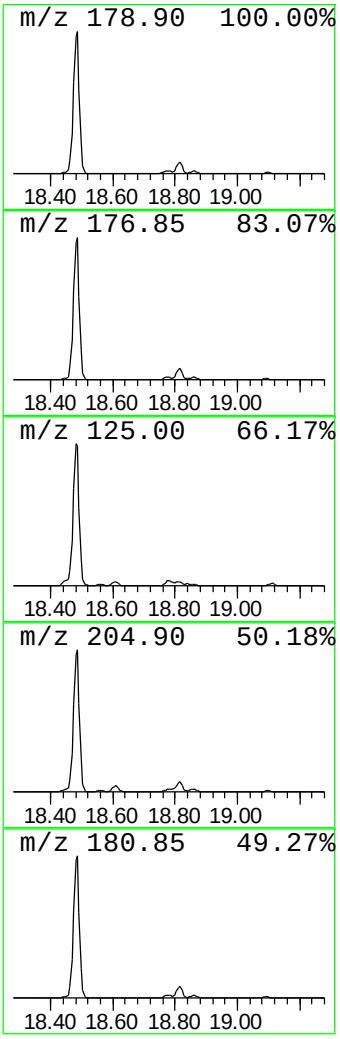
Peak Number 16 unknown-03 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.78	3.13 ng/ul	562469	Phenanthrene-d10	17.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	25
2	1H-Pyrrolo[3,4-c]pyridine-1,3(2H...	177	C8H7N3O2	090004-20-9	16
3	2,3-Dimethoxyphenylacetonitrile	177	C10H11NO2	004468-57-9	16
4	Benzohydrazide, 2-hydroxy-N2-[1-...	278	C14H18N2O4	1000258-69-9	12
5	Cyclopropane, pentachloro-	212	C3HCl5	006262-51-7	12





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Data File : BM016839.D
Acq On : 21 Sep 2018 19:17
Operator : SJ/JU
Sample : J5012-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08K5

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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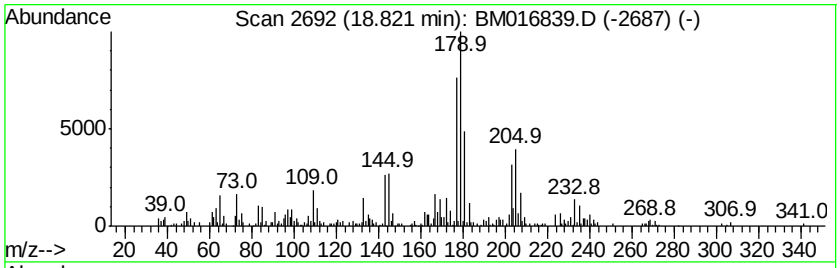
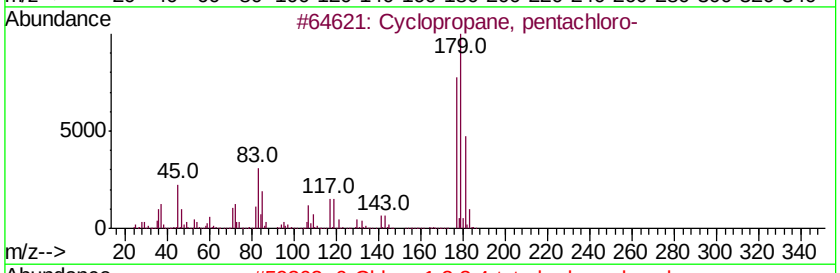
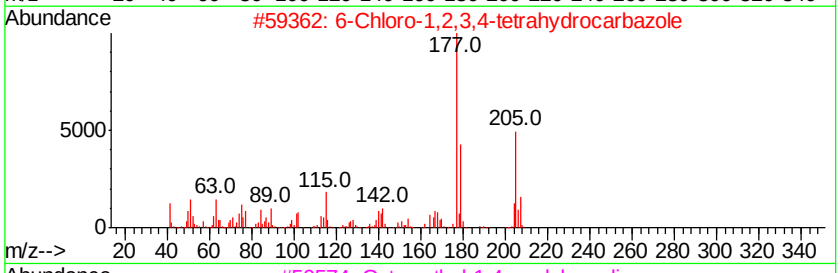
Peak Number 17 unknown-04 Concentration Rank 11

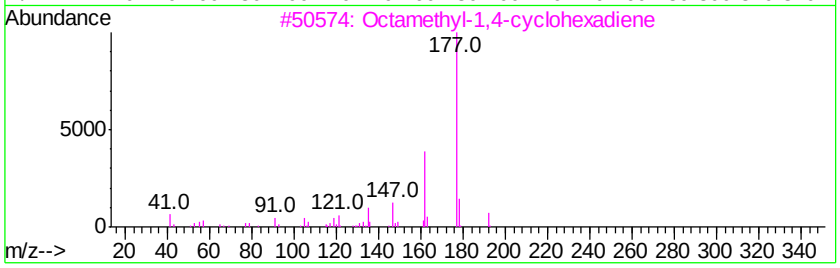
R.T.	EstConc	Area	Relative to ISTD	R.T.
18.82	7.33 ng/ul	1317480	Phenanthrene-d10	17.10

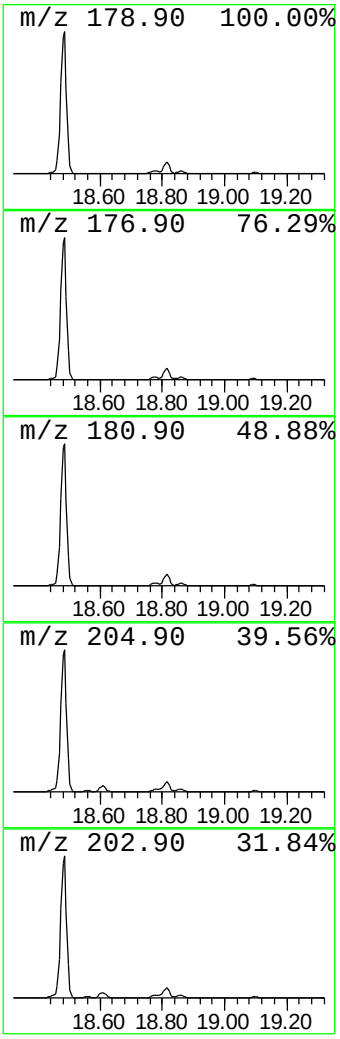
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, pentachloro-	212	C3HCl5	006262-51-7	38
2	6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	27
3	Octamethyl-1,4-cyclohexadiene	192	C14H24	172684-93-4	15
4	3-Buten-2-one, 4-(2,6,6-trimethy...	192	C13H20O	014901-07-6	15
5	5-Chloro-4-methyl-8-quinolinecar...	205	C11H8ClNO	028662-47-7	14











Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM092118\
Data File : BM016839.D
Acq On : 21 Sep 2018 19:17
Operator : SJ/JU
Sample : J5012-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08K5

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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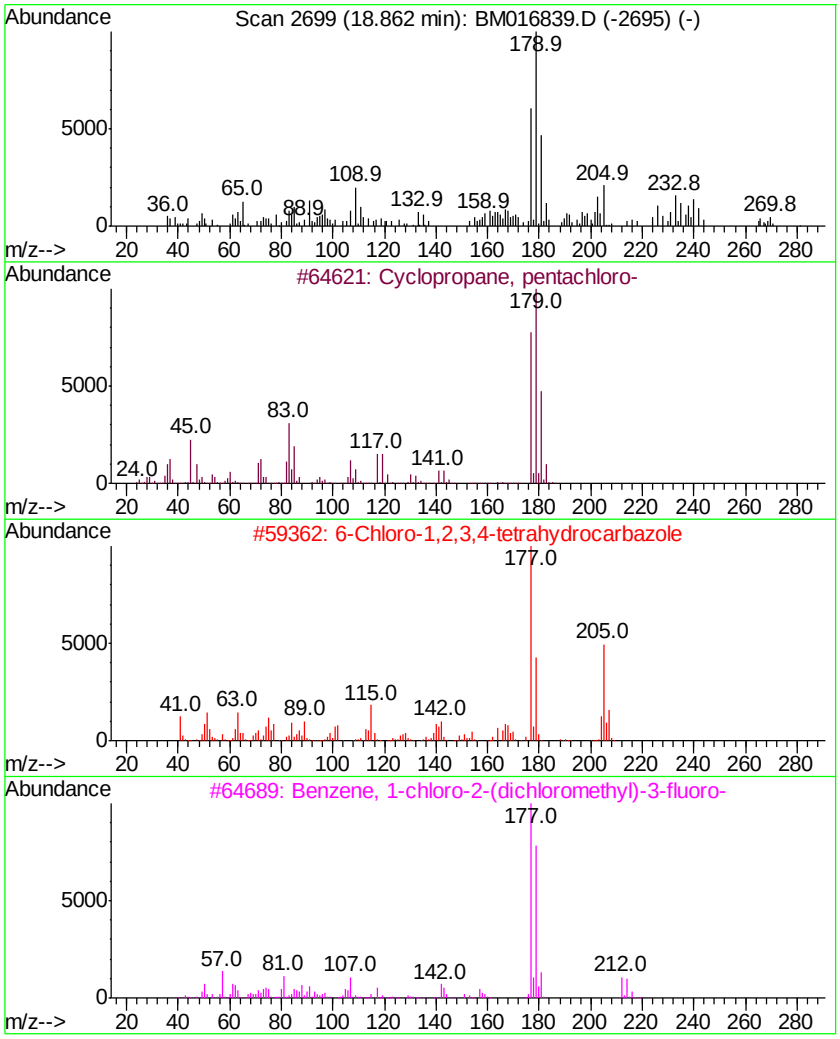
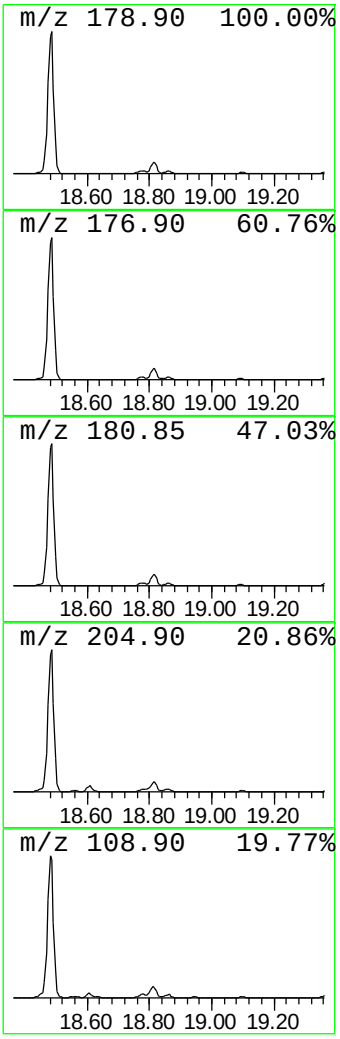
Peak Number 18 unknown-05 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.86	2.11 ng/ul	379983	Phenanthrene-d10	17.10

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, pentachloro-	212	C3HCl5	006262-51-7	40
2	6-Chloro-1,2,3,4-tetrahydrocarba...	205	C12H12ClN	036684-65-8	27
3	Benzene, 1-chloro-2-(dichloromet...	212	C7H4Cl3F	062476-62-4	25
4	6-Bromohexanoic acid, 4-hexadecy...	418	C22H43BrO2	1000282-90-6	16
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 : BM016839.D
Acq On : 21 Sep 2018 19:17
Operator : SJ/JU
Sample : J5012-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : C08K5

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, 1-bromo-...	9.22	15.9	ng/ul	1941030	2	10.52	2443540	20.0
Benzene, 1-bromo-...	9.53	8.3	ng/ul	1019410	2	10.52	2443540	20.0
Benzene, 1-chloro...	11.12	177.1	ng/ul	21642800	2	10.52	2443540	20.0
4-Chloro-2-nitrop...	11.80	6.4	ng/ul	776940	2	10.52	2443540	20.0
1(3H)-Isobenzofur...	12.76	2.1	ng/ul	332643	3	14.36	3190730	20.0
Benzene, 1,4-dich...	12.93	3.0	ng/ul	486021	3	14.36	3190730	20.0
unknown-01	18.49	98.0	ng/ul	17622300	4	17.10	3596680	20.0
unknown-02	18.60	4.1	ng/ul	732702	4	17.10	3596680	20.0
unknown-03	18.78	3.1	ng/ul	562469	4	17.10	3596680	20.0
unknown-04	18.82	7.3	ng/ul	1317480	4	17.10	3596680	20.0
unknown-05	18.86	2.1	ng/ul	379983	4	17.10	3596680	20.0