

Instrument :
BNA_M
ClientSampleId :
SSTD01002

mohammad
5/30/2017 4:40:19 PM

[illegible]

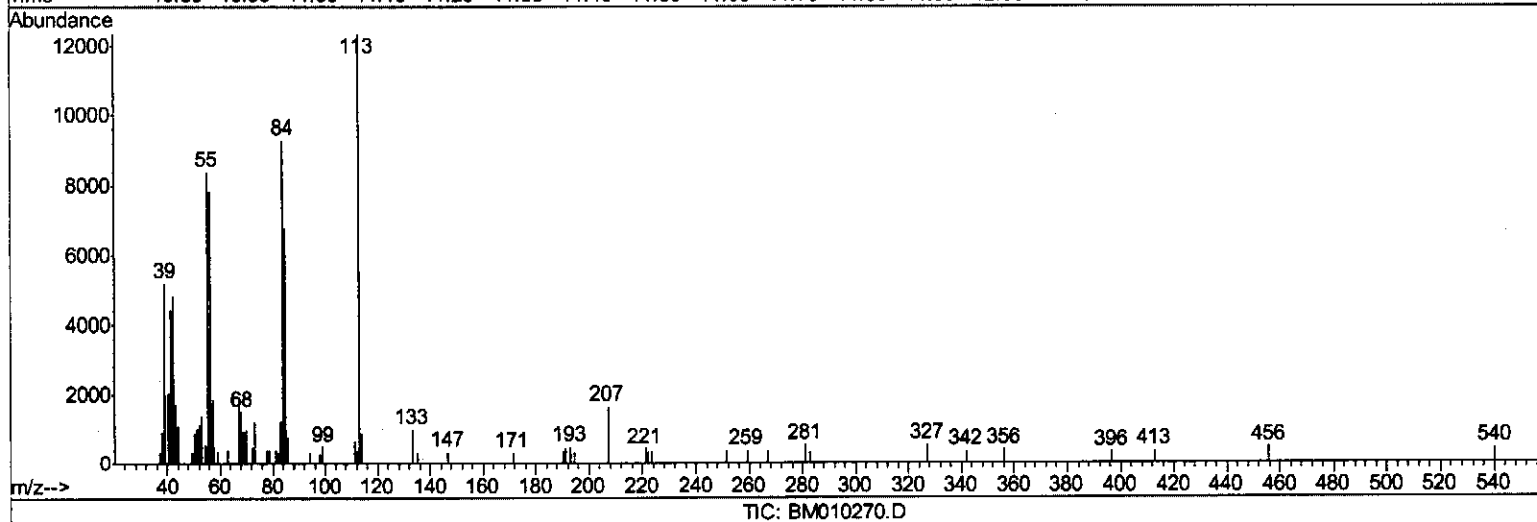
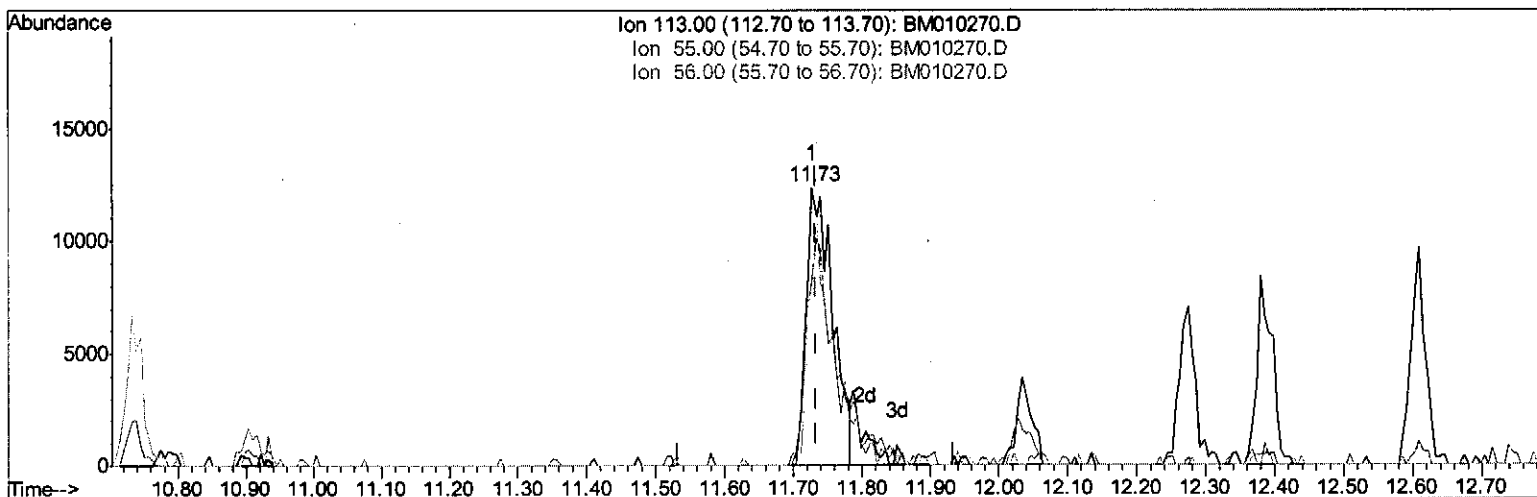
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052717\
Data File : BM010270.D
Acq On : 27 May 2017 11:54
Operator : SJ/MA
Sample : SSTD01002
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampled :
SSTD01002

Manual Integrations
APPROVED

mohammad
5/30/2017 4:40:19 PM

Quant Time: May 27 13:42:38 2017
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat May 27 13:30:54 2017
Response via : Initial Calibration



(32) Caprolactam

11.728min (-0.006) 8.87ng/ul

response 32804

Ion	Exp%	Act%
113.00	100	100
55.00	118.50	67.83#
56.00	107.50	63.31#
0.00	0.00	0.00

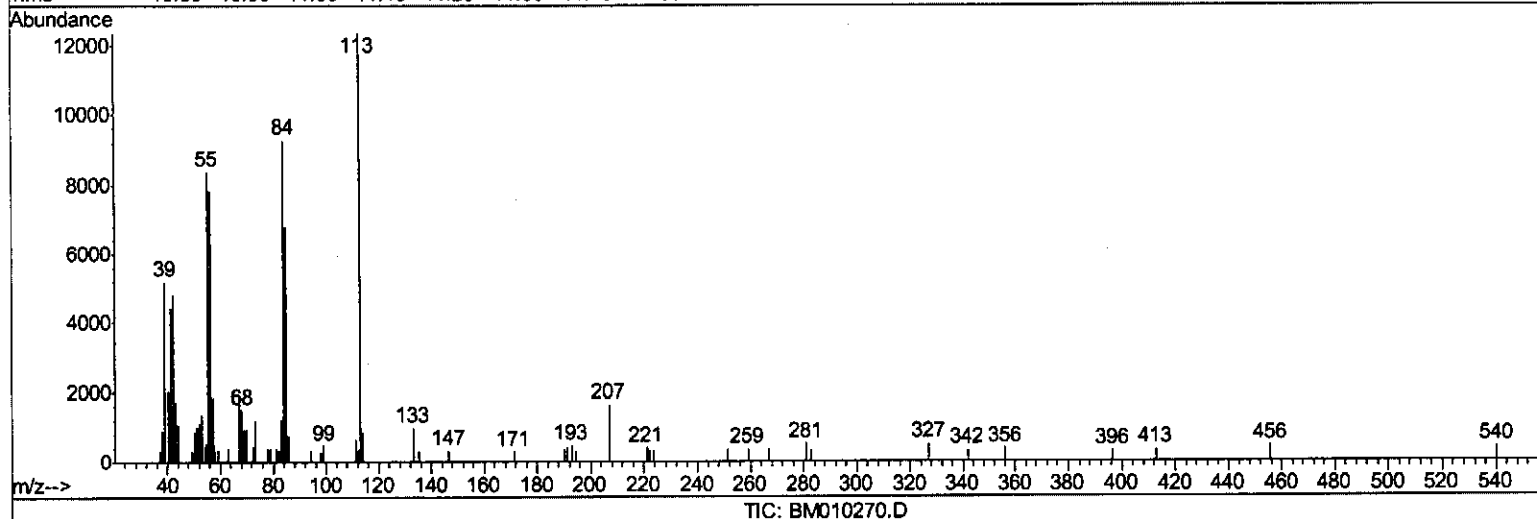
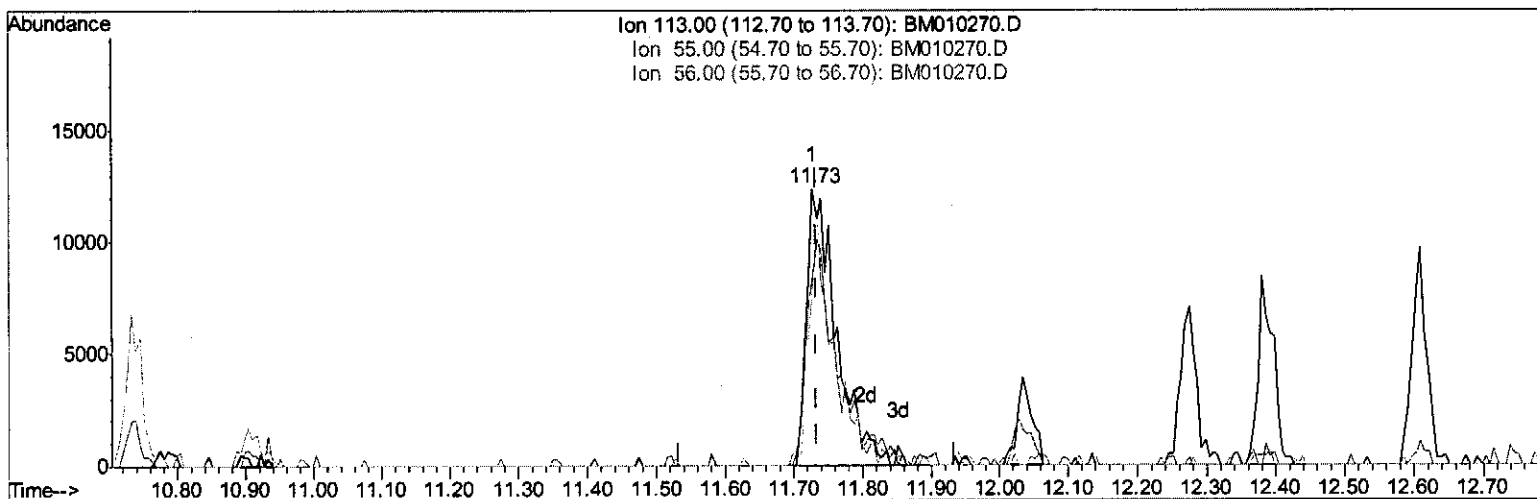
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052717\
Data File : BM010270.D
Acq On : 27 May 2017 11:54
Operator : SJ/MA
Sample : SSTD01002
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampled :
SSTD01002

Manual Integrations
APPROVED

mohammad
5/30/2017 4:40:19 PM

Quant Time: May 27 13:42:38 2017
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat May 27 13:30:54 2017
Response via : Initial Calibration



(32) Caprolactam

11.728min (-0.006) 10.00ng/ul m

SJ 5/30/17

response 36983

Ion	Exp%	Act%
113.00	100	100
55.00	118.50	67.83#
56.00	107.50	63.31#
0.00	0.00	0.00

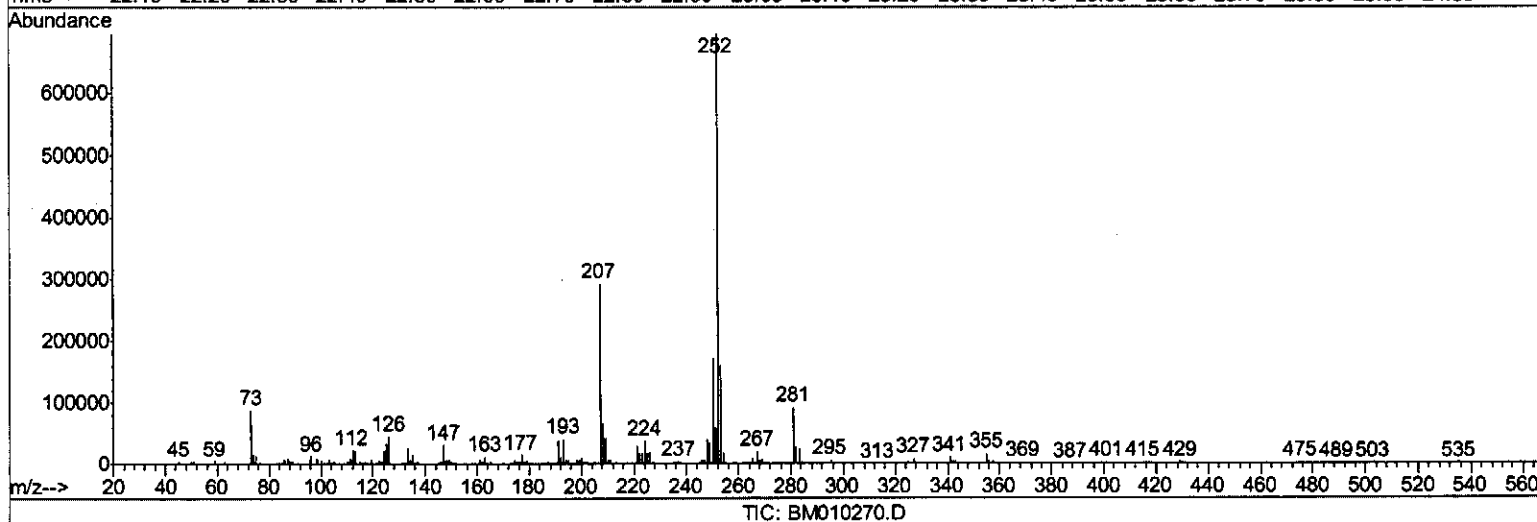
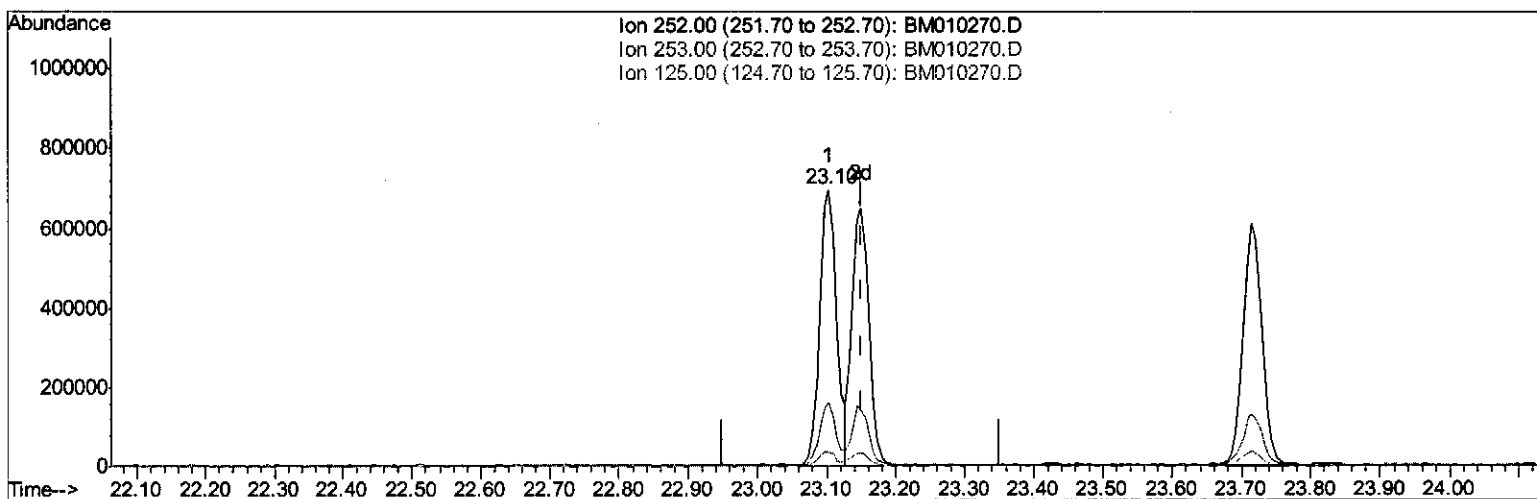
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052717\
Data File : BM010270.D
Acq On : 27 May 2017 11:54
Operator : SJ/MA
Sample : SSTD01002
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampled :
SSTD01002

Manual Integrations
APPROVED

mohammad
5/30/2017 4:40:19 PM

Quant Time: May 27 13:42:38 2017
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat May 27 13:30:54 2017
Response via : Initial Calibration



(86) Benzo(k)fluoranthene

23.103min (-0.047) 10.61ng/ul

response 1192828

Ion	Exp%	Act%
252.00	100	100
253.00	21.40	23.22
125.00	4.30	4.99
0.00	0.00	0.00

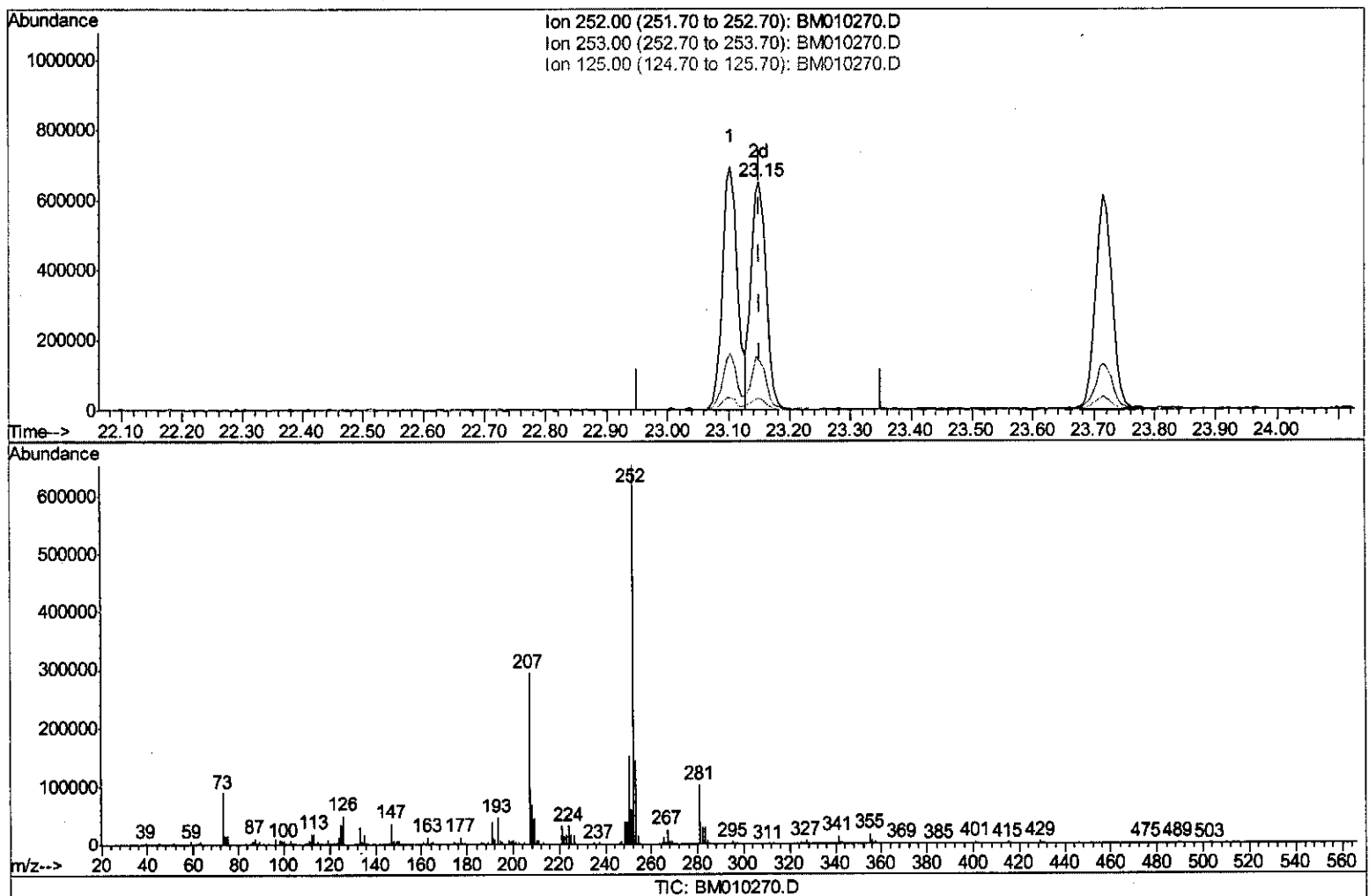
Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052717\
Data File : BM010270.D
Acq On : 27 May 2017 11:54
Operator : SJ/MA
Sample : SSTD01002
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampled :
SSTD01002

Manual Integrations
APPROVED

mohammad
5/30/2017 4:40:19 PM

Quant Time: May 27 13:42:38 2017
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat May 27 13:30:54 2017
Response via : Initial Calibration



(86) Benzo(k)fluoranthene

23.150min (+0.000) 9.90ng/ul m SJ 5/30/17

response 1112841

Ion	Exp%	Act%
252.00	100	100
253.00	21.40	21.97
125.00	4.30	5.23#
0.00	0.00	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052717\
 Data File : BM010270.D
 Acq On : 27 May 2017 11:54
 Operator : SJ/MA
 Sample : SSTD01002
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 Client Sampled :
 SSTD01002

Manual Integrations
 APPROVED

mohammad
 5/30/2017 4:40:19 PM

Quant Time: May 27 13:44:07 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat May 27 13:30:54 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	273646	20.00	ng/ul	0.00
18) Naphthalene-d8	10.74	136	997550	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.56	164	612348	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.31	188	1433786	20.00	ng/ul	0.00
75) Chrysene-d12	21.47	240	1993111	20.00	ng/ul	0.00
83) Perylene-d12	23.82	264	2142118	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.39	96	26703	4.10	ng/uL	0.00
5) Phenol-d5	7.12	99	182698	9.61	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.27	67	106536	9.73	ng/ul	0.00
9) 2-Chlorophenol-d4	7.47	132	152454	9.77	ng/ul	0.00
13) 4-Methylphenol-d8	8.65	113	150548	9.72	ng/ul	0.00
19) Nitrobenzene-d5	9.12	128	64232	9.29	ng/ul	0.00
22) 2-Nitrophenol-d4	9.83	143	75930	9.79	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.36	165	162606	9.87	ng/ul	0.00
29) 4-Chloroaniline-d4	10.90	131	133927	9.31	ng/ul	0.00
43) Dimethylphthalate-d6	13.98	166	482254	9.95	ng/ul	0.00
46) Acenaphthylene-d8	14.26	160	553509	9.89	ng/ul	0.00
51) 4-Nitrophenol-d4	14.80	143	57450	9.40	ng/ul	0.00
57) Fluorene-d10	15.56	176	429545	10.00	ng/ul	0.00
62) 4,6-Dinitro-2-methylphenol	15.68	200	79237	9.47	ng/ul	0.00
70) Anthracene-d10	17.41	188	652931	9.84	ng/ul	0.00
76) Pyrene-d10	19.69	212	776187	9.85	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.67	264	926343	9.76	ng/ul	0.00

Target Compounds

					Qvalue
2) 1,4-Dioxane	3.42	88	27524	4.05 ng/uL#	53
4) Benzaldehyde	7.10	77	132698	9.54 ng/ul	92
6) Phenol	7.15	94	188348	9.65 ng/ul	97
8) Bis(2-Chloroethyl)ether	7.36	93	133184	9.74 ng/ul	93
10) 2-Chlorophenol	7.50	128	152293	9.67 ng/ul	92
11) 2-Methylphenol	8.39	108	134851	9.55 ng/ul	95
12) 2,2'-oxybis(1-Chloropropan	8.46	45	60279	9.65 ng/ul#	64
14) Acetophenone	8.77	105	234228	9.50 ng/ul#	92
15) N-Nitroso-di-n-propylamine	8.74	70	120734	9.43 ng/ul#	83
16) 4-Methylphenol	8.72	108	150692	9.70 ng/ul	95
17) Hexachloroethane	9.00	117	70763	9.90 ng/ul	89
20) Nitrobenzene	9.16	77	190708	9.69 ng/ul	94
21) Isophorone	9.66	82	289150	9.62 ng/ul#	93
23) 2-Nitrophenol	9.86	139	83163	9.96 ng/ul	84
24) 2,4-Dimethylphenol	9.91	107	199231	9.86 ng/ul	96
25) Bis(2-Chloroethoxy)methane	10.15	93	172816	9.98 ng/ul	97
27) 2,4-Dichlorophenol	10.39	162	161322	9.98 ng/ul	94
28) Naphthalene	10.79	128	481980	9.89 ng/ul	97
30) 4-Chloroaniline	10.93	127	129806	9.19 ng/ul#	79
31) Hexachlorobutadiene	11.03	225	154616	9.98 ng/ul	92
32) Caprolactam	11.73	113	36983m	10.00 ng/ul	94
33) 4-Chloro-3-methylphenol	12.04	107	151592	9.95 ng/ul	94
34) 2-Methylnaphthalene	12.39	142	357908	9.86 ng/ul	93

SJ 5/30/17

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052717\
 Data File : BM010270.D
 Acq On : 27 May 2017 11:54
 Operator : SJ/MA
 Sample : SST01002
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleID :
 SST01002

Manual Integrations
 APPROVED

mohammad
 5/30/2017 4:40:19 PM

Quant Time: May 27 13:44:07 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM052717.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat May 27 13:30:54 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
36) 1,2,4,5-Tetrachlorobenzene	12.74	216	245018	9.88	ng/ul#	89
37) Hexachlorocyclopentadiene	12.70	237	151966	9.61	ng/ul#	97
38) 2,4,6-Trichlorophenol	13.00	196	131378	9.68	ng/ul	95
39) 2,4,5-Trichlorophenol	13.07	196	137267	9.90	ng/ul	91
40) 1,1'-Biphenyl	13.40	154	470675	9.90	ng/ul	95
41) 2-Chloronaphthalene	13.45	162	375077	9.96	ng/ul	98
42) 2-Nitroaniline	13.69	65	87365	9.95	ng/ul	96
44) Dimethylphthalate	14.03	163	473500	9.84	ng/ul	99
45) 2,6-Dinitrotoluene	14.16	165	80749	9.71	ng/ul#	82
47) Acenaphthylene	14.29	152	560153	10.25	ng/ul	97
48) 3-Nitroaniline	14.52	138	69872	9.28	ng/ul	84
49) Acenaphthene	14.63	153	391491	10.07	ng/ul	99
50) 2,4-Dinitrophenol	14.70	184	56002	9.87	ng/ul	97
52) 4-Nitrophenol	14.82	109	75531	9.36	ng/ul	88
53) Dibenzofuran	14.96	168	573059	9.97	ng/ul	93
54) 2,4-Dinitrotoluene	14.95	165	121407	10.00	ng/ul	95
55) 2,3,4,6-Tetrachlorophenol	15.19	232	129268	9.91	ng/ul#	94
56) Diethylphthalate	15.37	149	453307	9.90	ng/ul	96
58) Fluorene	15.61	166	463059	9.93	ng/ul	97
59) 4-Chlorophenyl-phenylether	15.60	204	268686	9.98	ng/ul	97
60) 4-Nitroaniline	15.68	138	73768	9.90	ng/ul	90
63) 4,6-Dinitro-2-methylphenol	15.69	198	86204	9.48	ng/ul#	95
64) N-Nitrosodiphenylamine	15.82	169	383270	9.57	ng/ul	98
65) 4-Bromophenyl-phenylether	16.49	248	162219	9.59	ng/ul#	89
66) Hexachlorobenzene	16.59	284	171427	9.58	ng/ul#	90
67) Atrazine	16.77	200	159743	9.88	ng/ul	92
68) Pentachlorophenol	16.95	266	92955	9.39	ng/ul	96
69) Phenanthrene	17.35	178	725625	9.85	ng/ul	99
71) Anthracene	17.44	178	748737	9.87	ng/ul	98
72) Carbazole	17.73	167	615614	9.87	ng/ul	99
73) Di-n-butylphthalate	18.24	149	685548	9.88	ng/ul	98
74) Fluoranthene	19.36	202	905359	9.97	ng/ul	98
77) Pyrene	19.72	202	955788	9.73	ng/ul	97
78) Butylbenzylphthalate	20.60	149	314635	9.84	ng/ul	97
79) 3,3'-Dichlorobenzidine	21.40	252	357138	9.69	ng/ul	95
80) Benzo(a)anthracene	21.46	228	1078832	9.87	ng/ul	98
81) Bis(2-ethylhexyl)phthalate	21.34	149	462828	9.77	ng/ul	100
82) Chrysene	21.51	228	976167	9.88	ng/ul	97
84) Di-n-octyl phthalate	22.24	149	859451	9.35	ng/ul#	93
85) Benzo(b)fluoranthene	23.10	252	1192828	9.94	ng/ul	98
86) Benzo(k)fluoranthene	23.15	252	1112841m)	9.90	ng/ul	97
88) Benzo(a)pyrene	23.72	252	1167342	9.89	ng/ul#	97
89) Indeno(1,2,3-cd)pyrene	26.21	276	1459324	9.79	ng/ul	98
90) Dibenzo(a,h)anthracene	26.23	278	1249522	9.76	ng/ul#	99
91) Benzo(g,h,i)perylene	26.97	276	1209381	9.74	ng/ul	100

(#) = qualifier out of range (m) = manual integration (+) = signals summed

SJ 5/30/17