

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031616\
 Data File : BM004629.D
 Acq On : 16 Mar 2016 10:48
 Operator : UM/SJ
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD02051

Quant Time: Mar 17 00:54:48 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM031516.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Mar 16 07:32:20 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.85	152	68866	20.00	ng/ul	0.00
18) Naphthalene-d8	10.63	136	291754	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.47	164	161633	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.22	188	357301	20.00	ng/ul	0.00
78) Chrysene-d12	21.40	240	415979	20.00	ng/ul	0.00
86) Perylene-d12	23.70	264	414848	20.00	ng/ul	0.00

System Monitoring Compounds

3) 1,4-Dioxane-d8	3.33	96	16133	8.08	ng/uL	0.00
5) Phenol-d5	7.00	99	116528	20.59	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.17	67	65949	20.56	ng/ul	0.00
9) 2-Chlorophenol-d4	7.37	132	92959	20.41	ng/ul	0.00
13) 4-Methylphenol-d8	8.54	113	89573	19.64	ng/ul	0.00
19) Nitrobenzene-d5	9.00	128	39199	19.32	ng/ul	0.00
22) 2-Nitrophenol-d4	9.72	143	42031	19.18	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.25	165	86281	20.07	ng/ul	0.00
29) 4-Chloroaniline-d4	10.77	131	116585	22.66	ng/ul	0.00
44) Dimethylphthalate-d6	13.89	166	245152	19.40	ng/ul	0.00
47) Acenaphthylene-d8	14.17	160	316219	20.41	ng/ul	0.00
52) 4-Nitrophenol-d4	14.66	143	44544	18.24	ng/ul	0.00
58) Fluorene-d10	15.47	176	220943	20.06	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.58	200	31622	16.84	ng/ul	0.00
71) Anthracene-d10	17.32	188	337756	20.53	ng/ul	0.00
79) Pyrene-d10	19.61	212	379474	19.98	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.55	264	373090	20.14	ng/ul	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.37	88	19140	8.22	ng/uL	92
4) Benzaldehyde	6.99	77	65144	20.76	ng/ul	99
6) Phenol	7.03	94	123644	20.77	ng/ul	98
8) Bis(2-Chloroethyl)ether	7.27	93	96381	20.80	ng/ul	96
10) 2-Chlorophenol	7.40	128	98099	20.40	ng/ul	95
11) 2-Methylphenol	8.27	108	90823	19.85	ng/ul	97
12) 2,2'-oxybis(1-Chloropropan	8.37	45	112663	20.32	ng/ul#	91
14) Acetophenone	8.66	105	146045	20.52	ng/ul	97
15) N-Nitroso-di-n-propylamine	8.65	70	68459	18.93	ng/ul	94
16) 4-Methylphenol	8.60	108	101325	20.11	ng/ul	98
17) Hexachloroethane	8.92	117	37198	19.47	ng/ul	97
20) Nitrobenzene	9.04	77	102052	20.38	ng/ul	96
21) Isophorone	9.56	82	185576	19.35	ng/ul	98
23) 2-Nitrophenol	9.75	139	48259	19.88	ng/ul	95
24) 2,4-Dimethylphenol	9.80	107	107570	20.23	ng/ul	99
25) Bis(2-Chloroethoxy)methane	10.05	93	123208	19.91	ng/ul	99
27) 2,4-Dichlorophenol	10.28	162	89493	20.21	ng/ul	99
28) Naphthalene	10.69	128	310256	20.39	ng/ul	98
30) 4-Chloroaniline	10.79	127	123610	22.87	ng/ul	98
31) Hexachlorobutadiene	10.97	225	54151	20.49	ng/ul	98
32) Caprolactam	11.55	113	26194	17.66	ng/ul	82
33) 4-Chloro-3-methylphenol	11.91	107	95075	18.97	ng/ul	99
34) 2-Methylnaphthalene	12.30	142	218302	20.15	ng/ul	100

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031616\
 Data File : BM004629.D
 Acq On : 16 Mar 2016 10:48
 Operator : UM/SJ
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTD02051

Quant Time: Mar 17 00:54:48 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM031516.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Mar 16 07:32:20 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
35) 1-Methylnaphthalene	12.52	142	209836	20.16	ng/ul	98
37) 1,2,4,5-Tetrachlorobenzene	12.66	216	107133	20.75	ng/ul	99
38) Hexachlorocyclopentadiene	12.65	237	51866	18.55	ng/ul	98
39) 2,4,6-Trichlorophenol	12.90	196	66861	19.61	ng/ul	100
40) 2,4,5-Trichlorophenol	12.97	196	72780	19.78	ng/ul	99
41) 1,1'-Biphenyl	13.31	154	276073	20.45	ng/ul	98
42) 2-Chloronaphthalene	13.35	162	208920	20.58	ng/ul	97
43) 2-Nitroaniline	13.55	65	50243	18.38	ng/ul	91
45) Dimethylphthalate	13.93	163	253580	19.77	ng/ul	99
46) 2,6-Dinitrotoluene	14.05	165	43912	19.23	ng/ul	89
48) Acenaphthylene	14.20	152	339513	20.40	ng/ul	100
49) 3-Nitroaniline	14.38	138	50567	20.50	ng/ul	93
50) Acenaphthene	14.54	153	228635	20.40	ng/ul	100
51) 2,4-Dinitrophenol	14.58	184	19579	15.06	ng/ul	91
53) 4-Nitrophenol	14.68	109	37665	19.02	ng/ul	93
54) Dibenzofuran	14.87	168	323670	20.29	ng/ul	96
55) 2,4-Dinitrotoluene	14.84	165	67017	19.46	ng/ul	94
56) 2,3,4,6-Tetrachlorophenol	15.10	232	62434	19.27	ng/ul	97
57) Diethylphthalate	15.30	149	251653	19.26	ng/ul	100
59) Fluorene	15.53	166	261026	20.49	ng/ul	98
60) 4-Chlorophenyl-phenylether	15.52	204	122828	20.22	ng/ul	93
61) 4-Nitroaniline	15.54	138	54694	20.61	ng/ul	95
64) 4,6-Dinitro-2-methylphenol	15.60	198	35577	17.44	ng/ul#	97
65) N-Nitrosodiphenylamine	15.73	169	221300	20.45	ng/ul	100
66) 4-Bromophenyl-phenylether	16.42	248	72942	20.27	ng/ul	93
67) Hexachlorobenzene	16.53	284	82324	20.38	ng/ul	94
68) Atrazine	16.68	200	75620	19.77	ng/ul	98
69) Pentachlorophenol	16.87	266	48755	18.61	ng/ul	98
70) Phenanthrene	17.26	178	418170	20.48	ng/ul	99
72) Anthracene	17.35	178	426829	20.76	ng/ul	99
73) 1,2,3,4-Tetrachlorobenzene	13.27	216	103569	21.07	ng/uL	99
74) Pentachlorobenzene	14.79	250	98544	20.77	ng/uL	99
75) Carbazole	17.62	167	391329	21.29	ng/ul	99
76) Di-n-butylphthalate	18.19	149	411058	15.77	ng/ul	100
77) Fluoranthene	19.27	202	478331	21.22	ng/ul	98
80) Pyrene	19.64	202	513604	20.37	ng/ul	98
81) Butylbenzylphthalate	20.54	149	183601	18.21	ng/ul	93
82) 3,3'-Dichlorobenzidine	21.32	252	144810	22.54	ng/ul	98
83) Benzo(a)anthracene	21.39	228	496604	20.45	ng/ul	100
84) Bis(2-ethylhexyl)phthalate	21.32	149	280066	19.54	ng/ul#	98
85) Chrysene	21.44	228	473052	20.40	ng/ul	100
87) Di-n-octyl phthalate	22.22	149	494622	19.95	ng/ul	97
88) Benzo(b)fluoranthene	23.01	252	482884	19.52	ng/ul	100
89) Benzo(k)fluoranthene	23.05	252	490089	20.84	ng/ul	99
91) Benzo(a)pyrene	23.60	252	479592	20.23	ng/ul	100
92) Indeno(1,2,3-cd)pyrene	26.03	276	533625	20.02	ng/ul	99
93) Dibenzo(a,h)anthracene	26.04	278	441736	20.03	ng/ul	99
94) Benzo(g,h,i)perylene	26.74	276	446022	19.76	ng/ul	99

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM031616\
Data File : BM004629.D
Acq On : 16 Mar 2016 10:48
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SSTD02051

Quant Time: Mar 17 00:54:48 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM031516.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Mar 16 07:32:20 2016
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)

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

