

Data Path : Z:\HPCHEM1\BNA G\DATA\BG100115\
 Data File : BG018981.D
 Acq On : 30 Sep 2015 20:41
 Operator : UM/NP
 Sample : SSTDC0020EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTD02021

Quant Time: Oct 01 05:41:29 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG091015.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 01 04:12:22 2015
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
 Max. RRF Dev : 20% Max. Rel. Area : 150%

	Compound	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	92	0.00
2	1,4-Dioxane	0.450	0.462	-2.7	91	0.00
3 S	1,4-Dioxane-d8	0.416	0.378	9.1	85	0.00
4	Benzaldehyde	0.950	1.065	-12.1	92	0.00
5 S	Phenol-d5	1.642	1.733	-5.5	98	0.00
6	Phenol	1.699	1.781	-4.8	99	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	0.952	0.924	2.9	88	0.00
8	Bis(2-Chloroethyl)ether	1.269	1.296	-2.1	93	0.00
9 S	2-Chlorophenol-d4	1.244	1.390	-11.7	104	0.00
10	2-Chlorophenol	1.288	1.392	-8.1	97	0.00
11	2-Methylphenol	1.306	1.382	-5.8	100	0.00
12	2,2'-oxybis(1-Chloropropane	1.506	1.402	6.9	85	0.00
13 S	4-Methylphenol-d8	1.333	1.418	-6.4	101	0.00
14	Acetophenone	2.030	2.200	-8.4	99	0.00
15 P	N-Nitroso-di-n-propylamine	1.047	1.000	4.5	88	0.00
16	4-Methylphenol	1.436	1.552	-8.1	100	0.00
17	Hexachloroethane	0.513	0.506	1.4	97	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	96	0.00
19 S	Nitrobenzene-d5	0.152	0.157	-3.3	102	0.00
20	Nitrobenzene	0.351	0.330	6.0	92	0.00
21	Isophorone	0.716	0.660	7.8	91	0.00
22 S	2-Nitrophenol-d4	0.152	0.175	-15.1	114	0.00
23 C	2-Nitrophenol	0.172	0.190	-10.5	112	0.00
24	2,4-Dimethylphenol	0.356	0.371	-4.2	105	0.00
25	Bis(2-Chloroethoxy)methane	0.422	0.413	2.1	96	0.00
26 S	2,4-Dichlorophenol-d3	0.321	0.344	-7.2	106	0.00
27 C	2,4-Dichlorophenol	0.321	0.342	-6.5	106	0.00
28	Naphthalene	1.018	1.029	-1.1	99	0.00
29 S	4-Chloroaniline-d4	0.373	0.460	-23.3#	114	0.00
30	4-Chloroaniline	0.386	0.464	-20.2#	110	0.00
31 C	Hexachlorobutadiene	0.198	0.212	-7.1	107	0.00
32	Caprolactam	0.131	0.124	5.3	90	0.01
33 C	4-Chloro-3-methylphenol	0.342	0.366	-7.0	105	0.00
34	2-Methylnaphthalene	0.752	0.780	-3.7	103	0.00
35	1-Methylnaphthalene	0.721	0.747	-3.6	104	0.00
36 I	Acenaphthene-d10	1.000	1.000	0.0	101	0.00
37	1,2,4,5-Tetrachlorobenzene	0.620	0.640	-3.2	110	0.00
38	Hexachlorocyclopentadiene	0.345	0.164	52.5#	52	0.00
39 C	2,4,6-Trichlorophenol	0.403	0.430	-6.7	108	0.00
40	2,4,5-Trichlorophenol	0.434	0.479	-10.4	113	0.00
41	1,1'-Biphenyl	1.514	1.520	-0.4	105	0.00
42	2-Chloronaphthalene	1.154	1.188	-2.9	106	0.00
43	2-Nitroaniline	0.342	0.328	4.1	101	0.00
44 S	Dimethylphthalate-d6	1.610	1.605	0.3	105	0.00

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Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
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	Compound	AvqRF	CCRF	%Dev	Area%	Dev(min)
45	Dimethylphthalate	1.560	1.558	0.1	104	0.00
46	2,6-Dinitrotoluene	0.314	0.360	-14.6	120	0.00
47 S	Acenaphthylene-d8	1.907	1.941	-1.8	105	0.00
48	Acenaphthylene	2.029	2.019	0.5	102	0.00
49	3-Nitroaniline	0.333	0.409	-22.8#	119	0.00
50 C	Acenaphthene	1.367	1.370	-0.2	104	0.00
51	2,4-Dinitrophenol	0.141	0.116	17.7	107	0.00
52 S	4-Nitrophenol-d4	0.309	0.274	11.3	93	0.00
53	4-Nitrophenol	0.212	0.176	17.0	87	0.00
54	Dibenzofuran	1.918	1.966	-2.5	106	0.00
55	2,4-Dinitrotoluene	0.467	0.535	-14.6	118	0.00
56	2,3,4,6-Tetrachlorophenol	0.419	0.415	1.0	105	0.00
57	Diethylphthalate	1.622	1.599	1.4	103	0.00
58 S	Fluorene-d10	1.407	1.464	-4.1	107	0.00
59	Fluorene	1.620	1.636	-1.0	104	0.00
60	4-Chlorophenyl-phenylether	0.772	0.795	-3.0	109	0.00
61	4-Nitroaniline	0.383	0.452	-18.0	113	0.00
62 I	Phenanthrene-d10	1.000	1.000	0.0	102	0.00
63 S	4,6-Dinitro-2-methylphenol-	0.095	0.098	-3.2	122	0.00
64	4,6-Dinitro-2-methylphenol	0.097	0.100	-3.1	119	0.00
65	N-Nitrosodiphenylamine	0.524	0.528	-0.8	107	0.00
66	4-Bromophenyl-phenylether	0.196	0.203	-3.6	113	0.00
67	Hexachlorobenzene	0.217	0.236	-8.8	120	0.00
68	Atrazine	0.231	0.237	-2.6	105	0.00
69 C	Pentachlorophenol	0.127	0.078	38.6#	68	0.00
70	Phenanthrene	1.017	1.025	-0.8	105	0.00
71 S	Anthracene-d10	0.881	0.899	-2.0	109	0.00
72	Anthracene	1.020	1.034	-1.4	105	0.00
73	1,2,3,4-Tetrachlorobenzene	0.238	0.240	-0.8	107	0.00
74	Pentachlorobenzene	0.232	0.259	-11.6	118	0.00
75	Carbazole	0.906	0.984	-8.6	104	0.00
76	Di-n-butylphthalate	1.144	1.165	-1.8	103	0.00
77 C	Fluoranthene	1.127	1.284	-13.9	105	0.00
78 I	Chrysene-d12	1.000	1.000	0.0	98	0.00
79 S	Pyrene-d10	0.920	0.993	-7.9	109	0.00
80	Pyrene	1.176	1.225	-4.2	102	0.00
81	Butylbenzylphthalate	0.451	0.489	-8.4	110	0.00
82	3,3'-Dichlorobenzidine	0.355	0.454	-27.9#	123	0.00
83	Benzo(a)anthracene	1.109	1.159	-4.5	104	0.00
84	Bis(2-ethylhexyl)phthalate	0.669	0.704	-5.2	105	0.00
85	Chrysene	1.023	1.065	-4.1	105	0.00
86 I	Perylene-d12	1.000	1.000	0.0	101	0.00
87	Di-n-octyl phthalate	1.058	1.163	-9.9	107	0.00

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88	Benzo(b)fluoranthene	1.131	1.173	-3.7	108	0.00
89	Benzo(k)fluoranthene	1.083	1.096	-1.2	102	0.00
90 S	Benzo(a)pyrene-d12	0.973	1.016	-4.4	109	0.00
91 C	Benzo(a)pyrene	1.075	1.110	-3.3	107	0.00
92	Indeno(1,2,3-cd)pyrene	1.246	1.301	-4.4	109	-0.02
93	Dibenzo(a,h)anthracene	1.000	1.069	-6.9	113	0.00
94	Benzo(a,h,i)perylene	1.040	1.075	-3.4	108	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 1