

Data Path : Z:\HPCHEM1\BNA_M\Data\BM042717\
 Data File : BM009731.D
 Acq On : 27 Apr 2017 09:03
 Operator : SJ/MA
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02032

Quant Time: Apr 27 18:09:06 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 27 18:07:34 2017
 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	112	0.00
2	1,4-Dioxane	0.535	0.514	3.9	102	0.00
3 S	1,4-Dioxane-d8	0.434	0.471	-8.5	117	0.00
4	Benzaldehyde	1.415	1.398	1.2	98	0.00
5 S	Phenol-d5	2.173	2.047	5.8	107	0.00
6	Phenol	2.246	2.181	2.9	111	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	1.482	1.451	2.1	109	0.00
8	Bis(2-Chloroethyl)ether	1.686	1.660	1.5	109	0.00
9 S	2-Chlorophenol-d4	1.382	1.382	0.0	112	0.00
10	2-Chlorophenol	1.415	1.449	-2.4	113	0.00
11	2-Methylphenol	1.616	1.557	3.7	109	0.00
12	2,2'-oxybis(1-Chloropropane	2.816	2.786	1.1	110	0.00
13 S	4-Methylphenol-d8	1.694	1.622	4.3	106	0.00
14	Acetophenone	2.786	2.682	3.7	107	0.00
15 P	N-Nitroso-di-n-propylamine	1.730	1.736	-0.3	111	0.00
16	4-Methylphenol	1.789	1.710	4.4	108	0.00
17	Hexachloroethane	0.657	0.674	-2.6	112	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	112	0.00
19 S	Nitrobenzene-d5	0.163	0.159	2.5	107	0.00
20	Nitrobenzene	0.529	0.513	3.0	109	0.00
21	Isophorone	0.936	0.907	3.1	107	0.00
22 S	2-Nitrophenol-d4	0.171	0.173	-1.2	114	0.00
23 C	2-Nitrophenol	0.187	0.180	3.7	108	0.00
24	2,4-Dimethylphenol	0.464	0.457	1.5	110	0.00
25	Bis(2-Chloroethoxy)methane	0.543	0.525	3.3	107	0.00
26 S	2,4-Dichlorophenol-d3	0.343	0.332	3.2	106	0.00
27 C	2,4-Dichlorophenol	0.338	0.324	4.1	109	0.00
28	Naphthalene	1.065	1.035	2.8	110	0.00
29 S	4-Chloroaniline-d4	0.413	0.429	-3.9	107	0.00
30	4-Chloroaniline	0.412	0.419	-1.7	105	0.00
31 C	Hexachlorobutadiene	0.236	0.228	3.4	111	0.00
32	Caprolactam	0.127	0.118	7.1	105	0.00
33 C	4-Chloro-3-methylphenol	0.421	0.412	2.1	108	0.00
34	2-Methylnaphthalene	0.819	0.789	3.7	107	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	108	0.00
36	1,2,4,5-Tetrachlorobenzene	0.678	0.667	1.6	109	0.00
37	Hexachlorocyclopentadiene	0.362	0.321	11.3	111	0.00
38 C	2,4,6-Trichlorophenol	0.461	0.449	2.6	108	0.00
39	2,4,5-Trichlorophenol	0.475	0.480	-1.1	111	0.00
40	1,1'-Biphenyl	1.645	1.592	3.2	107	0.00
41	2-Chloronaphthalene	1.265	1.234	2.5	107	0.00
42	2-Nitroaniline	0.539	0.539	0.0	107	0.00
43 S	Dimethylphthalate-d6	1.740	1.694	2.6	106	0.00
44	Dimethylphthalate	1.721	1.656	3.8	103	0.00

Data Path : Z:\HPCHEM1\BNA_M\Data\BM042717\
 Data File : BM009731.D
 Acq On : 27 Apr 2017 09:03
 Operator : SJ/MA
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02032

Quant Time: Apr 27 18:09:06 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 27 18:07:34 2017
 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	2,6-Dinitrotoluene	0.346	0.340	1.7	106	0.00
46 S	Acenaphthylene-d8	2.098	2.076	1.0	108	0.00
47	Acenaphthylene	2.069	2.000	3.3	105	0.00
48	3-Nitroaniline	0.341	0.318	6.7	95	0.00
49 C	Acenaphthene	1.407	1.386	1.5	108	0.00
50	2,4-Dinitrophenol	0.231	0.193	16.5	103	0.00
51 S	4-Nitrophenol-d4	0.336	0.321	4.5	108	0.00
52	4-Nitrophenol	0.385	0.367	4.7	111	0.00
53	Dibenzofuran	2.061	2.019	2.0	107	0.00
54	2,4-Dinitrotoluene	0.521	0.510	2.1	106	0.00
55	2,3,4,6-Tetrachlorophenol	0.461	0.450	2.4	106	0.00
56	Diethylphthalate	1.877	1.790	4.6	105	0.00
57 S	Fluorene-d10	1.570	1.501	4.4	105	0.00
58	Fluorene	1.759	1.719	2.3	109	0.00
59	4-Chlorophenyl-phenylether	0.911	0.874	4.1	105	0.00
60	4-Nitroaniline	0.398	0.331	16.8	90	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	108	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.141	0.121	14.2	101	0.00
63	4,6-Dinitro-2-methylphenol	0.146	0.135	7.5	106	0.00
64	N-Nitrosodiphenylamine	0.614	0.602	2.0	105	0.00
65	4-Bromophenyl-phenylether	0.228	0.226	0.9	106	0.00
66	Hexachlorobenzene	0.250	0.246	1.6	104	0.00
67	Atrazine	0.246	0.228	7.3	102	0.00
68 C	Pentachlorophenol	0.154	0.145	5.8	109	0.00
69	Phenanthrene	1.150	1.114	3.1	104	0.00
70 S	Anthracene-d10	1.027	1.006	2.0	105	0.00
71	Anthracene	1.203	1.178	2.1	107	0.00
72	Carbazole	1.082	1.022	5.5	104	0.00
73	Di-n-butylphthalate	1.318	1.277	3.1	104	0.00
74 C	Fluoranthene	1.450	1.376	5.1	105	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	105	0.00
76 S	Pyrene-d10	0.883	0.875	0.9	104	0.00
77	Pyrene	1.130	1.121	0.8	104	0.00
78	Butylbenzylphthalate	0.468	0.464	0.9	104	0.00
79	3,3'-Dichlorobenzidine	0.370	0.362	2.2	92	0.00
80	Benzo(a)anthracene	1.186	1.168	1.5	103	0.00
81	Bis(2-ethylhexyl)phthalate	0.710	0.681	4.1	101	0.00
82	Chrysene	1.068	1.027	3.8	100	0.00
83 I	Perylene-d12	1.000	1.000	0.0	96	0.00
84	Di-n-octyl phthalate	1.461	1.349	7.7	100	0.00
85	Benzo(b)fluoranthene	1.299	1.301	-0.2	101	0.00
86	Benzo(k)fluoranthene	1.190	1.174	1.3	96	0.00
87 S	Benzo(a)pyrene-d12	0.973	0.965	0.8	97	0.00

Data Path : Z:\HPCHEM1\BNA_M\Data\BM042717\
Data File : BM009731.D
Acq On : 27 Apr 2017 09:03
Operator : SJ/MA
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02032

Quant Time: Apr 27 18:09:06 2017
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 27 18:07:34 2017
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	AvgRF	CCRF	%Dev	Area%	Dev(min)
88 C	Benzo(a)pyrene	1.208	1.193	1.2	96	0.00
89	Indeno(1,2,3-cd)pyrene	1.308	1.161	11.2	84	0.00
90	Dibenzo(a,h)anthracene	1.114	1.013	9.1	86	0.00
91	Benzo(g,h,i)perylene	1.052	0.906	13.9	81	0.00

(#) = Out of Range

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