

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM042517\
 Data File : BM009713.D
 Acq On : 25 Apr 2017 18:34
 Operator : SJ/MA
 Sample : SSTDC0020EC
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02029

Quant Time: Apr 26 03:39:44 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 25 13:32:43 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	95	0.00
2	1,4-Dioxane	0.535	0.520	2.8	87	0.00
3 S	1,4-Dioxane-d8	0.434	0.453	-4.4	95	0.00
4	Benzaldehyde	1.415	0.985	30.4#	58	0.00
5 S	Phenol-d5	2.173	2.027	6.7	90	0.00
6	Phenol	2.246	2.171	3.3	93	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	1.482	1.459	1.6	93	0.00
8	Bis(2-Chloroethyl)ether	1.686	1.640	2.7	91	0.00
9 S	2-Chlorophenol-d4	1.382	1.377	0.4	94	0.00
10	2-Chlorophenol	1.415	1.389	1.8	92	0.00
11	2-Methylphenol	1.616	1.522	5.8	90	0.00
12	2,2'-oxybis(1-Chloropropane	2.816	2.789	1.0	93	0.00
13 S	4-Methylphenol-d8	1.694	1.635	3.5	90	0.00
14	Acetophenone	2.786	2.673	4.1	91	0.00
15 P	N-Nitroso-di-n-propylamine	1.730	1.705	1.4	92	0.00
16	4-Methylphenol	1.789	1.749	2.2	93	0.00
17	Hexachloroethane	0.657	0.656	0.2	92	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	93	0.00
19 S	Nitrobenzene-d5	0.163	0.155	4.9	87	0.00
20	Nitrobenzene	0.529	0.528	0.2	94	0.00
21	Isophorone	0.936	0.921	1.6	91	0.00
22 S	2-Nitrophenol-d4	0.171	0.169	1.2	93	0.00
23 C	2-Nitrophenol	0.187	0.179	4.3	90	0.00
24	2,4-Dimethylphenol	0.464	0.463	0.2	93	0.00
25	Bis(2-Chloroethoxy)methane	0.543	0.523	3.7	89	0.00
26 S	2,4-Dichlorophenol-d3	0.343	0.352	-2.6	94	0.00
27 C	2,4-Dichlorophenol	0.338	0.328	3.0	92	0.00
28	Naphthalene	1.065	1.036	2.7	92	0.00
29 S	4-Chloroaniline-d4	0.413	0.406	1.7	85	0.00
30	4-Chloroaniline	0.412	0.403	2.2	85	0.00
31 C	Hexachlorobutadiene	0.236	0.234	0.8	95	0.00
32	Caprolactam	0.127	0.120	5.5	90	0.00
33 C	4-Chloro-3-methylphenol	0.421	0.428	-1.7	93	0.00
34	2-Methylnaphthalene	0.819	0.786	4.0	89	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	92	0.00
36	1,2,4,5-Tetrachlorobenzene	0.678	0.689	-1.6	95	0.00
37	Hexachlorocyclopentadiene	0.362	0.322	11.0	94	0.00
38 C	2,4,6-Trichlorophenol	0.461	0.455	1.3	93	0.00
39	2,4,5-Trichlorophenol	0.475	0.465	2.1	91	0.00
40	1,1'-Biphenyl	1.645	1.612	2.0	92	0.00
41	2-Chloronaphthalene	1.265	1.262	0.2	93	0.00
42	2-Nitroaniline	0.539	0.550	-2.0	93	0.00
43 S	Dimethylphthalate-d6	1.740	1.720	1.1	92	0.00
44	Dimethylphthalate	1.721	1.676	2.6	89	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM042517\
 Data File : BM009713.D
 Acq On : 25 Apr 2017 18:34
 Operator : SJ/MA
 Sample : SSTDC0020EC
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02029

Quant Time: Apr 26 03:39:44 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Apr 25 13:32:43 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	0.347	-0.3	92	0.00
46 S	Acenaphthylene-d8	2.098	2.101	-0.1	93	0.00
47	Acenaphthylene	2.069	2.058	0.5	92	0.00
48	3-Nitroaniline	0.341	0.302	11.4	77	0.00
49 C	Acenaphthene	1.407	1.411	-0.3	93	0.00
50	2,4-Dinitrophenol	0.231	0.200	13.4	90	0.00
51 S	4-Nitrophenol-d4	0.336	0.327	2.7	93	0.00
52	4-Nitrophenol	0.385	0.382	0.8	98	0.00
53	Dibenzofuran	2.061	2.056	0.2	92	0.00
54	2,4-Dinitrotoluene	0.521	0.519	0.4	92	0.00
55	2,3,4,6-Tetrachlorophenol	0.461	0.467	-1.3	94	0.00
56	Diethylphthalate	1.877	1.816	3.2	91	0.00
57 S	Fluorene-d10	1.570	1.554	1.0	93	0.00
58	Fluorene	1.759	1.745	0.8	94	0.00
59	4-Chlorophenyl-phenylether	0.911	0.894	1.9	91	0.00
60	4-Nitroaniline	0.398	0.310	22.1#	72	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	96	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.141	0.126	10.6	94	0.00
63	4,6-Dinitro-2-methylphenol	0.146	0.132	9.6	93	0.00
64	N-Nitrosodiphenylamine	0.614	0.592	3.6	93	0.00
65	4-Bromophenyl-phenylether	0.228	0.222	2.6	94	0.00
66	Hexachlorobenzene	0.250	0.242	3.2	92	0.00
67	Atrazine	0.246	0.235	4.5	94	0.00
68 C	Pentachlorophenol	0.154	0.139	9.7	93	0.00
69	Phenanthrene	1.150	1.112	3.3	93	0.00
70 S	Anthracene-d10	1.027	1.013	1.4	94	0.00
71	Anthracene	1.203	1.176	2.2	95	0.00
72	Carbazole	1.082	1.030	4.8	94	0.00
73	Di-n-butylphthalate	1.318	1.276	3.2	93	0.00
74 C	Fluoranthene	1.450	1.394	3.9	95	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	97	0.00
76 S	Pyrene-d10	0.883	0.863	2.3	95	0.00
77	Pyrene	1.130	1.106	2.1	96	0.00
78	Butylbenzylphthalate	0.468	0.456	2.6	96	0.00
79	3,3'-Dichlorobenzidine	0.370	0.335	9.5	79	0.00
80	Benzo(a)anthracene	1.186	1.175	0.9	96	0.00
81	Bis(2-ethylhexyl)phthalate	0.710	0.680	4.2	94	0.00
82	Chrysene	1.068	1.036	3.0	94	0.00
83 I	Perylene-d12	1.000	1.000	0.0	95	0.00
84	Di-n-octyl phthalate	1.461	1.269	13.1	93	0.00
85	Benzo(b)fluoranthene	1.299	1.269	2.3	97	0.00
86	Benzo(k)fluoranthene	1.190	1.189	0.1	96	0.00
87 S	Benzo(a)pyrene-d12	0.973	0.966	0.7	96	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM042517\
Data File : BM009713.D
Acq On : 25 Apr 2017 18:34
Operator : SJ/MA
Sample : SSTDCCC020EC
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02029

Quant Time: Apr 26 03:39:44 2017
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM042517.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Apr 25 13:32:43 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	1.201	0.6	95	0.00
89	Indeno(1,2,3-cd)pyrene	1.308	1.312	-0.3	94	0.00
90	Dibenzo(a,h)anthracene	1.114	1.110	0.4	93	0.00
91	Benzo(g,h,i)perylene	1.052	1.054	-0.2	93	0.00

(#) = Out of Range

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