

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM091417\
 Data File : BM011537.D
 Acq On : 14 Sep 2017 12:42
 Operator : SJ/JU
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02017

Quant Time: Sep 14 17:30:00 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM090817MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 13 07:51:23 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	94	0.00
2	1,4-Dioxane	0.487	0.495	-1.6	95	0.00
3 S	1,4-Dioxane-d8	0.445	0.453	-1.8	95	0.00
4	Benzaldehyde	1.097	1.054	3.9	84	0.00
5 S	Phenol-d5	1.920	1.907	0.7	97	0.00
6	Phenol	1.906	1.917	-0.6	97	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	1.294	1.362	-5.3	99	0.00
8	Bis(2-Chloroethyl)ether	1.501	1.527	-1.7	97	0.00
9 S	2-Chlorophenol-d4	1.447	1.466	-1.3	95	0.00
10	2-Chlorophenol	1.412	1.431	-1.3	96	0.00
11	2-Methylphenol	1.445	1.443	0.1	97	0.00
12	2,2'-oxybis(1-Chloropropane	2.433	2.674	-9.9	103	0.00
13 S	4-Methylphenol-d8	1.499	1.496	0.2	96	0.00
14	Acetophenone	2.292	2.323	-1.4	97	0.00
15 P	N-Nitroso-di-n-propylamine	1.230	1.306	-6.2	99	0.00
16	4-Methylphenol	1.582	1.598	-1.0	97	0.00
17	Hexachloroethane	0.540	0.546	-1.1	97	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	96	0.00
19 S	Nitrobenzene-d5	0.152	0.152	0.0	96	0.00
20	Nitrobenzene	0.357	0.372	-4.2	100	0.00
21	Isophorone	0.719	0.745	-3.6	99	0.00
22 S	2-Nitrophenol-d4	0.158	0.154	2.5	95	0.00
23 C	2-Nitrophenol	0.166	0.165	0.6	95	0.00
24	2,4-Dimethylphenol	0.355	0.361	-1.7	97	0.00
25	Bis(2-Chloroethoxy)methane	0.469	0.468	0.2	96	0.00
26 S	2,4-Dichlorophenol-d3	0.318	0.316	0.6	95	0.00
27 C	2,4-Dichlorophenol	0.306	0.306	0.0	95	0.00
28	Naphthalene	1.037	1.034	0.3	96	0.00
29 S	4-Chloroaniline-d4	0.351	0.407	-16.0	126	0.00
30	4-Chloroaniline	0.333	0.388	-16.5	125	0.00
31 C	Hexachlorobutadiene	0.174	0.176	-1.1	99	0.00
32	Caprolactam	0.096	0.093	3.1	96	0.00
33 C	4-Chloro-3-methylphenol	0.325	0.333	-2.5	99	0.00
34	2-Methylnaphthalene	0.770	0.771	-0.1	96	0.00
35	1-Methylnaphthalene	0.734	0.732	0.3	96	0.00
36 I	Acenaphthene-d10	1.000	1.000	0.0	98	0.00
37	1,2,4,5-Tetrachlorobenzene	0.593	0.604	-1.9	97	0.00
38	Hexachlorocyclopentadiene	0.333	0.314	5.7	97	0.00
39 C	2,4,6-Trichlorophenol	0.404	0.398	1.5	95	0.00
40	2,4,5-Trichlorophenol	0.410	0.413	-0.7	97	0.00
41	1,1'-Biphenyl	1.664	1.657	0.4	96	0.00
42	2-Chloronaphthalene	1.252	1.244	0.6	96	0.00
43	2-Nitroaniline	0.389	0.419	-7.7	102	0.00
44 S	Dimethylphthalate-d6	1.691	1.673	1.1	95	0.00

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM091417\
 Data File : BM011537.D
 Acq On : 14 Sep 2017 12:42
 Operator : SJ/JU
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02017

Quant Time: Sep 14 17:30:00 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM090817MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 13 07:51:23 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
45	Dimethylphthalate	1.619	1.598	1.3	94	0.00
46	2,6-Dinitrotoluene	0.286	0.287	-0.3	96	0.00
47 S	Acenaphthylene-d8	2.045	2.064	-0.9	96	0.00
48	Acenaphthylene	2.058	2.081	-1.1	96	0.00
49	3-Nitroaniline	0.353	0.332	5.9	95	0.00
50 C	Acenaphthene	1.391	1.383	0.6	96	0.00
51	2,4-Dinitrophenol	0.177	0.149	15.8	94	0.00
52 S	4-Nitrophenol-d4	0.318	0.293	7.9	94	0.00
53	4-Nitrophenol	0.204	0.201	1.5	101	0.00
54	Dibenzofuran	1.945	1.948	-0.2	96	0.00
55	2,4-Dinitrotoluene	0.420	0.421	-0.2	96	0.00
56	2,3,4,6-Tetrachlorophenol	0.373	0.367	1.6	95	0.00
57	Diethylphthalate	1.685	1.663	1.3	95	0.00
58 S	Fluorene-d10	1.465	1.469	-0.3	97	0.00
59	Fluorene	1.630	1.679	-3.0	98	0.00
60	4-Chlorophenyl-phenylether	0.764	0.779	-2.0	98	0.00
61	4-Nitroaniline	0.378	0.340	10.1	89	0.00
62 I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
63 S	4,6-Dinitro-2-methylphenol-	0.116	0.101	12.9	94	0.00
64	4,6-Dinitro-2-methylphenol	0.119	0.106	10.9	93	0.00
65	N-Nitrosodiphenylamine	0.609	0.608	0.2	96	0.00
66	4-Bromophenyl-phenylether	0.204	0.203	0.5	97	0.00
67	Hexachlorobenzene	0.225	0.221	1.8	97	0.00
68	Atrazine	0.211	0.207	1.9	98	0.00
69 C	Pentachlorophenol	0.144	0.130	9.7	96	0.00
70	Phenanthrene	1.115	1.129	-1.3	96	0.00
71 S	Anthracene-d10	0.985	0.991	-0.6	97	0.00
72	Anthracene	1.144	1.168	-2.1	96	0.00
73	1,2,3,4-Tetrachlorobenzene	0.267	0.265	0.7	96	0.00
74	Pentachlorobenzene	0.270	0.269	0.4	96	0.00
75	Carbazole	0.976	1.026	-5.1	96	0.00
76	Di-n-butylphthalate	1.282	1.347	-5.1	95	0.00
77 C	Fluoranthene	1.147	1.291	-12.6	97	0.00
78 I	Chrysene-d12	1.000	1.000	0.0	97	0.00
79 S	Pyrene-d10	0.936	0.962	-2.8	98	0.00
80	Pyrene	1.167	1.221	-4.6	96	0.00
81	Butylbenzylphthalate	0.537	0.542	-0.9	97	0.00
82	3,3'-Dichlorobenzidine	0.363	0.343	5.5	90	0.00
83	Benzo(a)anthracene	1.101	1.140	-3.5	96	0.00
84	Bis(2-ethylhexyl)phthalate	0.824	0.844	-2.4	95	0.00
85	Chrysene	0.998	1.016	-1.8	95	0.00
86 I	Perylene-d12	1.000	1.000	0.0	85	0.00
87	Di-n-octyl phthalate	1.419	1.738	-22.5#	94	0.00

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA M\DATA\BM091417\
 Data File : BM011537.D
 Acq On : 14 Sep 2017 12:42
 Operator : SJ/JU
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02017

Quant Time: Sep 14 17:30:00 2017
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-BM090817MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 13 07:51:23 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	Benzo(b)fluoranthene	1.203	1.281	-6.5	91	0.00
89	Benzo(k)fluoranthene	1.156	1.174	-1.6	85	0.00
90 S	Benzo(a)pyrene-d12	0.953	0.967	-1.5	86	0.00
91 C	Benzo(a)pyrene	1.136	1.151	-1.3	85	0.00
92	Indeno(1,2,3-cd)pyrene	1.249	1.129	9.6	75	0.00
93	Dibenzo(a,h)anthracene	1.059	0.960	9.3	76	0.00
94	Benzo(a,h,i)perylene	1.012	0.896	11.5	74	0.00

(#) = Out of Range

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