

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062017\
 Data File : BM010601.D
 Acq On : 20 Jun 2017 21:55
 Operator : SJ/MA
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02038

Quant Time: Jun 21 01:31:55 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jun 21 01:19:58 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	74	0.00
2	1,4-Dioxane	0.390	0.392	-0.5	67	0.00
3 S	1,4-Dioxane-d8	0.349	0.345	1.1	68	0.00
4	Benzaldehyde	1.121	1.212	-8.1	66	0.00
5 S	Phenol-d5	1.905	1.748	8.2	68	0.00
6	Phenol	1.962	1.776	9.5	68	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	1.088	0.991	8.9	64	0.00
8	Bis(2-Chloroethyl)ether	1.429	1.311	8.3	67	0.00
9 S	2-Chlorophenol-d4	1.337	1.301	2.7	71	0.00
10	2-Chlorophenol	1.336	1.303	2.5	70	0.00
11	2-Methylphenol	1.496	1.419	5.1	69	0.00
12	2,2'-oxybis(1-Chloropropane	0.997	0.906	9.1	64	0.00
13 S	4-Methylphenol-d8	1.580	1.502	4.9	69	0.00
14	Acetophenone	2.579	2.457	4.7	70	0.00
15 P	N-Nitroso-di-n-propylamine	1.390	1.324	4.7	68	0.00
16	4-Methylphenol	1.646	1.558	5.3	69	0.00
17	Hexachloroethane	0.594	0.581	2.2	71	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	75	0.00
19 S	Nitrobenzene-d5	0.143	0.142	0.7	75	0.00
20	Nitrobenzene	0.416	0.398	4.3	70	0.00
21	Isophorone	0.828	0.751	9.3	65	0.00
22 S	2-Nitrophenol-d4	0.164	0.166	-1.2	74	0.00
23 C	2-Nitrophenol	0.174	0.173	0.6	71	0.00
24	2,4-Dimethylphenol	0.443	0.431	2.7	72	0.00
25	Bis(2-Chloroethoxy)methane	0.459	0.422	8.1	66	0.00
26 S	2,4-Dichlorophenol-d3	0.346	0.341	1.4	73	0.00
27 C	2,4-Dichlorophenol	0.337	0.331	1.8	72	0.00
28	Naphthalene	0.995	0.977	1.8	72	0.00
29 S	4-Chloroaniline-d4	0.365	0.406	-11.2	71	0.00
30	4-Chloroaniline	0.367	0.403	-9.8	71	0.00
31 C	Hexachlorobutadiene	0.254	0.262	-3.1	75	0.00
32	Caprolactam	0.118	0.114	3.4	68	0.00
33 C	4-Chloro-3-methylphenol	0.430	0.418	2.8	70	0.00
34	2-Methylnaphthalene	0.800	0.794	0.8	73	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	75	0.00
36	1,2,4,5-Tetrachlorobenzene	0.706	0.705	0.1	74	0.00
37	Hexachlorocyclopentadiene	0.406	0.401	1.2	79	0.00
38 C	2,4,6-Trichlorophenol	0.482	0.486	-0.8	75	0.00
39	2,4,5-Trichlorophenol	0.498	0.480	3.6	72	0.00
40	1,1'-Biphenyl	1.564	1.541	1.5	74	0.00
41	2-Chloronaphthalene	1.230	1.190	3.3	73	0.00
42	2-Nitroaniline	0.360	0.343	4.7	71	0.00
43 S	Dimethylphthalate-d6	1.756	1.755	0.1	74	0.00
44	Dimethylphthalate	1.721	1.713	0.5	74	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062017\
 Data File : BM010601.D
 Acq On : 20 Jun 2017 21:55
 Operator : SJ/MA
 Sample : SSTDC0020
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02038

Quant Time: Jun 21 01:31:55 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jun 21 01:19:58 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	0.298	0.314	-5.4	80	0.00
46 S	Acenaphthylene-d8	2.042	2.021	1.0	74	0.00
47	Acenaphthylene	1.945	1.913	1.6	74	0.00
48	3-Nitroaniline	0.301	0.292	3.0	72	0.00
49 C	Acenaphthene	1.312	1.276	2.7	73	0.00
50	2,4-Dinitrophenol	0.216	0.200	7.4	80	0.00
51 S	4-Nitrophenol-d4	0.309	0.289	6.5	73	0.00
52	4-Nitrophenol	0.392	0.410	-4.6	78	0.00
53	Dibenzofuran	1.987	1.958	1.5	73	0.00
54	2,4-Dinitrotoluene	0.462	0.473	-2.4	76	0.00
55	2,3,4,6-Tetrachlorophenol	0.504	0.511	-1.4	75	0.00
56	Diethylphthalate	1.811	1.769	2.3	73	0.00
57 S	Fluorene-d10	1.518	1.525	-0.5	75	0.00
58	Fluorene	1.664	1.614	3.0	73	0.00
59	4-Chlorophenyl-phenylether	0.904	0.901	0.3	75	0.00
60	4-Nitroaniline	0.355	0.348	2.0	76	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	76	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.123	0.113	8.1	76	0.00
63	4,6-Dinitro-2-methylphenol	0.131	0.124	5.3	78	0.00
64	N-Nitrosodiphenylamine	0.555	0.547	1.4	74	0.00
65	4-Bromophenyl-phenylether	0.229	0.224	2.2	75	0.00
66	Hexachlorobenzene	0.241	0.241	0.0	77	0.00
67	Atrazine	0.234	0.238	-1.7	76	0.00
68 C	Pentachlorophenol	0.170	0.165	2.9	75	0.00
69	Phenanthrene	1.063	1.063	0.0	76	0.00
70 S	Anthracene-d10	0.966	0.961	0.5	74	0.00
71	Anthracene	1.095	1.099	-0.4	75	0.00
72	Carbazole	0.956	0.931	2.6	73	0.00
73	Di-n-butylphthalate	1.215	1.202	1.1	74	0.00
74 C	Fluoranthene	1.370	1.361	0.7	75	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	80	0.00
76 S	Pyrene-d10	0.929	0.919	1.1	76	0.00
77	Pyrene	1.155	1.136	1.6	76	0.00
78	Butylbenzylphthalate	0.434	0.435	-0.2	78	0.00
79	3,3'-Dichlorobenzidine	0.400	0.403	-0.8	77	0.00
80	Benzo(a)anthracene	1.165	1.171	-0.5	78	0.00
81	Bis(2-ethylhexyl)phthalate	0.663	0.640	3.5	75	0.00
82	Chrysene	1.038	1.052	-1.3	78	0.00
83 I	Perylene-d12	1.000	1.000	0.0	81	0.00
84	Di-n-octyl phthalate	1.398	1.195	14.5	76	0.00
85	Benzo(b)fluoranthene	1.298	1.290	0.6	81	0.00
86	Benzo(k)fluoranthene	1.231	1.226	0.4	80	0.00
87 S	Benzo(a)pyrene-d12	0.967	0.968	-0.1	81	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062017\
Data File : BM010601.D
Acq On : 20 Jun 2017 21:55
Operator : SJ/MA
Sample : SSTDCCC020
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02038

Quant Time: Jun 21 01:31:55 2017
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jun 21 01:19:58 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	1.195	1.205	-0.8	81	0.00
89	Indeno(1,2,3-cd)pyrene	1.240	1.286	-3.7	82	0.00
90	Dibenzo(a,h)anthracene	1.033	1.063	-2.9	81	0.00
91	Benzo(g,h,i)perylene	0.987	1.024	-3.7	82	0.00

(#) = Out of Range

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