

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_M\DATA\BM062617\
 Data File : BM010761.D
 Acq On : 27 Jun 2017 00:39
 Operator : SJ/JU
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02002

Quant Time: Jun 27 02:26:46 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 27 01:28:19 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	92	0.00
2	1,4-Dioxane	0.390	0.381	2.3	82	0.00
3 S	1,4-Dioxane-d8	0.349	0.356	-2.0	88	0.00
4	Benzaldehyde	1.121	1.240	-10.6	84	0.00
5 S	Phenol-d5	1.905	1.696	11.0	82	0.00
6	Phenol	1.962	1.765	10.0	84	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	1.088	0.952	12.5	77	0.00
8	Bis(2-Chloroethyl)ether	1.429	1.288	9.9	82	0.00
9 S	2-Chlorophenol-d4	1.337	1.304	2.5	89	0.00
10	2-Chlorophenol	1.336	1.282	4.0	86	0.00
11	2-Methylphenol	1.496	1.404	6.1	86	0.00
12	2,2'-oxybis(1-Chloropropane	0.997	0.770	22.8#	68	0.00
13 S	4-Methylphenol-d8	1.580	1.444	8.6	83	0.00
14	Acetophenone	2.579	2.403	6.8	85	0.00
15 P	N-Nitroso-di-n-propylamine	1.390	1.288	7.3	82	0.00
16	4-Methylphenol	1.646	1.528	7.2	85	0.00
17	Hexachloroethane	0.594	0.599	-0.8	91	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	91	0.00
19 S	Nitrobenzene-d5	0.143	0.145	-1.4	93	0.00
20	Nitrobenzene	0.416	0.427	-2.6	91	0.00
21	Isophorone	0.828	0.742	10.4	79	0.00
22 S	2-Nitrophenol-d4	0.164	0.176	-7.3	96	0.00
23 C	2-Nitrophenol	0.174	0.179	-2.9	89	0.00
24	2,4-Dimethylphenol	0.443	0.449	-1.4	91	0.00
25	Bis(2-Chloroethoxy)methane	0.459	0.415	9.6	80	0.00
26 S	2,4-Dichlorophenol-d3	0.346	0.362	-4.6	94	0.00
27 C	2,4-Dichlorophenol	0.337	0.341	-1.2	91	0.00
28	Naphthalene	0.995	0.974	2.1	88	0.00
29 S	4-Chloroaniline-d4	0.365	0.410	-12.3	87	0.00
30	4-Chloroaniline	0.367	0.404	-10.1	87	0.00
31 C	Hexachlorobutadiene	0.254	0.274	-7.9	96	0.00
32	Caprolactam	0.118	0.108	8.5	78	0.00
33 C	4-Chloro-3-methylphenol	0.430	0.416	3.3	84	0.00
34	2-Methylnaphthalene	0.800	0.786	1.8	88	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	91	0.00
36	1,2,4,5-Tetrachlorobenzene	0.706	0.743	-5.2	94	0.00
37	Hexachlorocyclopentadiene	0.406	0.346	14.8	82	0.00
38 C	2,4,6-Trichlorophenol	0.482	0.496	-2.9	93	0.00
39	2,4,5-Trichlorophenol	0.498	0.510	-2.4	92	0.00
40	1,1'-Biphenyl	1.564	1.541	1.5	90	0.00
41	2-Chloronaphthalene	1.230	1.220	0.8	90	0.00
42	2-Nitroaniline	0.360	0.387	-7.5	98	0.00
43 S	Dimethylphthalate-d6	1.756	1.719	2.1	88	0.00
44	Dimethylphthalate	1.721	1.675	2.7	87	0.00

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_M\DATA\BM062617\
 Data File : BM010761.D
 Acq On : 27 Jun 2017 00:39
 Operator : SJ/JU
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02002

Quant Time: Jun 27 02:26:46 2017
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 27 01:28:19 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	0.339	-13.8	104	0.00
46 S	Acenaphthylene-d8	2.042	2.022	1.0	90	0.00
47	Acenaphthylene	1.945	1.907	2.0	89	0.00
48	3-Nitroaniline	0.301	0.332	-10.3	99	0.00
49 C	Acenaphthene	1.312	1.277	2.7	89	0.00
50	2,4-Dinitrophenol	0.216	0.222	-2.8	107	0.00
51 S	4-Nitrophenol-d4	0.309	0.309	0.0	94	0.00
52	4-Nitrophenol	0.392	0.463	-18.1	107	0.00
53	Dibenzofuran	1.987	1.987	0.0	90	0.00
54	2,4-Dinitrotoluene	0.462	0.500	-8.2	98	0.00
55	2,3,4,6-Tetrachlorophenol	0.504	0.511	-1.4	91	0.00
56	Diethylphthalate	1.811	1.744	3.7	87	0.00
57 S	Fluorene-d10	1.518	1.523	-0.3	91	0.00
58	Fluorene	1.664	1.646	1.1	90	0.00
59	4-Chlorophenyl-phenylether	0.904	0.910	-0.7	92	0.00
60	4-Nitroaniline	0.355	0.366	-3.1	97	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	95	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.123	0.123	0.0	104	0.00
63	4,6-Dinitro-2-methylphenol	0.131	0.129	1.5	102	0.00
64	N-Nitrosodiphenylamine	0.555	0.523	5.8	89	0.00
65	4-Bromophenyl-phenylether	0.229	0.225	1.7	94	0.00
66	Hexachlorobenzene	0.241	0.230	4.6	91	0.00
67	Atrazine	0.234	0.237	-1.3	94	0.00
68 C	Pentachlorophenol	0.170	0.159	6.5	90	0.00
69	Phenanthrene	1.063	1.034	2.7	92	0.00
70 S	Anthracene-d10	0.966	0.950	1.7	92	0.00
71	Anthracene	1.095	1.057	3.5	90	0.00
72	Carbazole	0.956	0.908	5.0	89	0.00
73	Di-n-butylphthalate	1.215	1.139	6.3	87	0.00
74 C	Fluoranthene	1.370	1.356	1.0	94	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	93	0.00
76 S	Pyrene-d10	0.929	0.953	-2.6	91	0.00
77	Pyrene	1.155	1.176	-1.8	91	0.00
78	Butylbenzylphthalate	0.434	0.465	-7.1	95	0.00
79	3,3'-Dichlorobenzidine	0.400	0.407	-1.7	90	0.00
80	Benzo(a)anthracene	1.165	1.192	-2.3	92	0.00
81	Bis(2-ethylhexyl)phthalate	0.663	0.674	-1.7	91	0.00
82	Chrysene	1.038	1.051	-1.3	90	0.00
83 I	Perylene-d12	1.000	1.000	0.0	77	0.00
84	Di-n-octyl phthalate	1.398	1.520	-8.7	92	0.00
85	Benzo(b)fluoranthene	1.298	1.366	-5.2	82	0.00
86	Benzo(k)fluoranthene	1.231	1.265	-2.8	79	0.00
87 S	Benzo(a)pyrene-d12	0.967	0.982	-1.6	78	0.00

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_M\DATA\BM062617\
Data File : BM010761.D
Acq On : 27 Jun 2017 00:39
Operator : SJ/JU
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02002

Quant Time: Jun 27 02:26:46 2017
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM062017.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jun 27 01:28:19 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	1.207	-1.0	78	0.00
89	Indeno(1,2,3-cd)pyrene	1.240	1.182	4.7	72	0.00
90	Dibenzo(a,h)anthracene	1.033	0.989	4.3	72	0.00
91	Benzo(g,h,i)perylene	0.987	0.966	2.1	74	0.00

(#) = Out of Range

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