

Data Path : S:\HPCHEM1\BNA_M\DATA\BM072015\
 Data File : BM001995.D
 Acq On : 21 Jul 2015 04:07
 Operator : TP/UM
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02059

Quant Time: Jul 21 04:49:04 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM071315.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 21 01:37:44 2015
 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	135	0.00
2	1,4-Dioxane	8.000	6.860	14.2	118	0.00
3 S	1,4-Dioxane-d8	8.000	7.083	11.5	118	0.00
4	Benzaldehyde	20.000	16.652	16.7	133	0.00
5 S	Phenol-d5	20.000	18.514	7.4	130	0.00
6	Phenol	20.000	18.534	7.3	130	0.01
7 S	Bis-(2-Chloroethyl)ether-d8	20.000	18.246	8.8	123	0.00
8	Bis(2-Chloroethyl)ether	20.000	19.000	5.0	128	0.00
9 S	2-Chlorophenol-d4	20.000	19.329	3.4	133	0.00
10	2-Chlorophenol	20.000	19.467	2.7	134	0.00
11	2-Methylphenol	20.000	18.648	6.8	129	0.01
12	2,2'-oxybis(1-Chloropropane	20.000	16.939	15.3	114	0.00
13 S	4-Methylphenol-d8	20.000	18.573	7.1	128	0.00
14	Acetophenone	20.000	18.377	8.1	126	0.00
15 P	N-Nitroso-di-n-propylamine	20.000	17.629	11.9	118	0.00
16	4-Methylphenol	20.000	18.714	6.4	129	0.00
17	Hexachloroethane	20.000	17.767	11.2	123	0.00
18 I	Naphthalene-d8	20.000	20.000	0.0	130	0.00
19 S	Nitrobenzene-d5	20.000	19.723	1.4	131	0.00
20	Nitrobenzene	20.000	18.789	6.1	124	0.00
21	Isophorone	20.000	17.928	10.4	118	0.00
22 S	2-Nitrophenol-d4	20.000	18.417	7.9	122	0.00
23 C	2-Nitrophenol	20.000	17.883	10.6	117	0.00
24	2,4-Dimethylphenol	20.000	18.555	7.2	123	0.00
25	Bis(2-Chloroethoxy)methane	20.000	18.533	7.3	124	0.00
26 S	2,4-Dichlorophenol-d3	20.000	19.715	1.4	130	0.00
27 C	2,4-Dichlorophenol	20.000	19.275	3.6	128	0.00
28	Naphthalene	20.000	19.149	4.3	129	0.00
29 S	4-Chloroaniline-d4	20.000	20.827	-4.1	132	0.00
30	4-Chloroaniline	20.000	20.217	-1.1	129	0.00
31 C	Hexachlorobutadiene	20.000	19.736	1.3	134	0.00
32	Caprolactam	20.000	17.519	12.4	115	0.02
33 C	4-Chloro-3-methylphenol	20.000	17.678	11.6	118	0.01
34	2-Methylnaphthalene	20.000	18.731	6.3	125	0.00
35 I	Acenaphthene-d10	20.000	20.000	0.0	118	0.00
36	1,2,4,5-Tetrachlorobenzene	20.000	20.991	-5.0	129	0.00
37	Hexachlorocyclopentadiene	20.000	6.858	65.7#	43	0.00
38 C	2,4,6-Trichlorophenol	20.000	19.910	0.4	121	0.00
39	2,4,5-Trichlorophenol	20.000	19.042	4.8	115	0.01
40	1,1'-Biphenyl	20.000	20.144	-0.7	123	0.00
41	2-Chloronaphthalene	20.000	20.152	-0.8	123	0.01
42	2-Nitroaniline	20.000	18.311	8.4	109	0.00
43 S	Dimethylphthalate-d6	20.000	18.811	5.9	114	0.00
44	Dimethylphthalate	20.000	18.703	6.5	113	0.00

Data Path : S:\HPCHEM1\BNA_M\DATA\BM072015\
 Data File : BM001995.D
 Acq On : 21 Jul 2015 04:07
 Operator : TP/UM
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02059

Quant Time: Jul 21 04:49:04 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM071315.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 21 01:37:44 2015
 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	Amount	Calc.	%Dev	Area%	Dev(min)
45	2,6-Dinitrotoluene	20.000	18.779	6.1	112	0.00
46 S	Acenaphthylene-d8	20.000	19.488	2.6	119	0.00
47	Acenaphthylene	20.000	19.404	3.0	118	0.00
48	3-Nitroaniline	20.000	21.031	-5.2	130	0.01
49 C	Acenaphthene	20.000	19.257	3.7	118	0.00
50	2,4-Dinitrophenol	20.000	2.178	89.1#	14	0.01
51 S	4-Nitrophenol-d4	20.000	13.808	31.0#	83	0.01
52	4-Nitrophenol	20.000	13.275	33.6#	80	0.01
53	Dibenzofuran	20.000	18.975	5.1	116	0.00
54	2,4-Dinitrotoluene	20.000	17.505	12.5	105	0.00
55	2,3,4,6-Tetrachlorophenol	20.000	16.676	16.6	100	0.00
56	Diethylphthalate	20.000	18.318	8.4	111	0.00
57 S	Fluorene-d10	20.000	18.806	6.0	114	0.00
58	Fluorene	20.000	18.634	6.8	114	0.00
59	4-Chlorophenyl-phenylether	20.000	19.118	4.4	118	0.00
60	4-Nitroaniline	20.000	19.472	2.6	122	0.00
61 I	Phenanthrene-d10	20.000	20.000	0.0	104	0.00
62 S	4,6-Dinitro-2-methylphenol-	20.000	3.863	80.7#	21	0.00
63	4,6-Dinitro-2-methylphenol	20.000	3.849	80.8#	21	0.00
64	N-Nitrosodiphenylamine	20.000	20.799	-4.0	112	0.00
65	4-Bromophenyl-phenylether	20.000	21.063	-5.3	112	0.00
66	Hexachlorobenzene	20.000	20.367	-1.8	111	0.00
67	Atrazine	20.000	20.483	-2.4	107	0.00
68 C	Pentachlorophenol	20.000	13.619	31.9#	75	0.00
69	Phenanthrene	20.000	19.077	4.6	103	0.00
70 S	Anthracene-d10	20.000	19.177	4.1	104	0.00
71	Anthracene	20.000	19.073	4.6	103	0.00
72	1,2,3,4-Tetrachlorobenzene	20.000	23.397	-17.0	128	0.00
73	Pentachlorobenzene	20.000	22.147	-10.7	120	0.00
74	Carbazole	20.000	18.472	7.6	98	0.00
75	Di-n-butylphthalate	20.000	20.047	-0.2	106	0.00
76 C	Fluoranthene	20.000	16.675	16.6	88	0.00
77 I	Chrysene-d12	20.000	20.000	0.0	73	0.00
78 S	Pyrene-d10	20.000	22.136	-10.7	84	0.00
79	Pyrene	20.000	21.957	-9.8	83	0.00
80	Butylbenzylphthalate	20.000	22.900	-14.5	85	0.00
81	3,3'-Dichlorobenzidine	20.000	16.497	17.5	64	0.00
82	Benzo(a)anthracene	20.000	19.437	2.8	74	0.00
83	Bis(2-ethylhexyl)phthalate	20.000	22.002	-10.0	84	0.00
84	Chrysene	20.000	19.204	4.0	72	0.00
85 I	Perylene-d12	20.000	20.000	0.0	79	0.00
86	Di-n-octyl phthalate	20.000	19.208	4.0	79	0.00
87	Benzo(b)fluoranthene	20.000	18.258	8.7	74	0.00

Data Path : S:\HPCHEM1\BNA_M\DATA\BM072015\
Data File : BM001995.D
Acq On : 21 Jul 2015 04:07
Operator : TP/UM
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02059

Quant Time: Jul 21 04:49:04 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM071315.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jul 21 01:37:44 2015
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	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	20.000	19.134	4.3	76	0.00
89 S	Benzo(a)pyrene-d12	20.000	19.344	3.3	78	0.00
90 C	Benzo(a)pyrene	20.000	19.090	4.6	77	0.00
91	Indeno(1,2,3-cd)pyrene	20.000	21.533	-7.7	91	0.00
92	Dibenzo(a,h)anthracene	20.000	21.788	-8.9	92	0.01
93	Benzo(g,h,i)perylene	20.000	22.224	-11.1	95	0.01

(#) = Out of Range

SPCC's out = 0 CCC's out = 1