

Data Path : Z:\HPCHEM1\BNA M\DATA\BM042015\
 Data File : BM001099.D
 Acq On : 21 Apr 2015 11:45
 Operator : TP/IZ
 Sample : SSTD02042
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Apr 21 19:13:04 2015
 Quant Method : Z:\HPCHEM1\BNA M\Methods\SOM01.2-EPA-BM041515.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 16 05:47:45 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	85	-0.01
2	1,4-Dioxane	0.444	0.428	3.6	84	-0.01
3 S	1,4-Dioxane-d8	0.429	0.433	-0.9	88	0.00
4	Benzaldehyde	1.001	1.004	-0.3	83	-0.01
5 S	Phenol-d5	1.560	1.569	-0.6	84	0.00
6	Phenol	1.650	1.647	0.2	85	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	0.972	0.942	3.1	82	0.00
8	Bis(2-Chloroethyl)ether	1.316	1.291	1.9	84	-0.01
9 S	2-Chlorophenol-d4	1.269	1.322	-4.2	88	-0.01
10	2-Chlorophenol	1.343	1.369	-1.9	88	0.00
11	2-Methylphenol	1.245	1.267	-1.8	85	0.00
12	2,2'-oxybis(1-Chloropropane	2.345	2.120	9.6	76	-0.01
13 S	4-Methylphenol-d8	1.231	1.256	-2.0	85	0.00
14	Acetophenone	1.997	1.958	2.0	82	-0.01
15 P	N-Nitroso-di-n-propylamine	0.993	0.961	3.2	81	-0.01
16	4-Methylphenol	1.358	1.351	0.5	84	0.00
17	Hexachloroethane	0.521	0.526	-1.0	87	-0.01
18 I	Naphthalene-d8	1.000	1.000	0.0	83	-0.01
19 S	Nitrobenzene-d5	0.135	0.143	-5.9	87	-0.01
20	Nitrobenzene	0.349	0.345	1.1	81	-0.01
21	Isophorone	0.616	0.619	-0.5	81	-0.01
22 S	2-Nitrophenol-d4	0.151	0.163	-7.9	87	-0.01
23 C	2-Nitrophenol	0.166	0.175	-5.4	87	0.00
24	2,4-Dimethylphenol	0.350	0.347	0.9	82	0.00
25	Bis(2-Chloroethoxy)methane	0.407	0.397	2.5	80	-0.01
26 S	2,4-Dichlorophenol-d3	0.294	0.307	-4.4	86	0.00
27 C	2,4-Dichlorophenol	0.293	0.298	-1.7	84	-0.01
28	Naphthalene	1.037	1.028	0.9	82	-0.01
29 S	4-Chloroaniline-d4	0.286	0.389	-36.0#	113	-0.01
30	4-Chloroaniline	0.291	0.391	-34.4#	111	-0.01
31 C	Hexachlorobutadiene	0.176	0.175	0.6	82	-0.01
32	Caprolactam	0.078	0.090	-15.4	96	-0.01
33 C	4-Chloro-3-methylphenol	0.301	0.310	-3.0	85	0.00
34	2-Methylnaphthalene	0.724	0.703	2.9	81	-0.01
35 I	Acenaphthene-d10	1.000	1.000	0.0	82	-0.01
36	1,2,4,5-Tetrachlorobenzene	0.617	0.616	0.2	82	-0.01
37	Hexachlorocyclopentadiene	0.323	0.242	25.1#	59	-0.01
38 C	2,4,6-Trichlorophenol	0.377	0.410	-8.8	88	0.00
39	2,4,5-Trichlorophenol	0.405	0.439	-8.4	86	0.00
40	1,1'-Biphenyl	1.675	1.654	1.3	80	-0.01
41	2-Chloronaphthalene	1.281	1.279	0.2	81	-0.01
42	2-Nitroaniline	0.348	0.364	-4.6	89	0.00
43 S	Dimethylphthalate-d6	1.333	1.348	-1.1	84	0.00
44	Dimethylphthalate	1.493	1.493	0.0	83	0.00

Data Path : Z:\HPCHEM1\BNA M\DATA\BM042015\
 Data File : BM001099.D
 Acq On : 21 Apr 2015 11:45
 Operator : TP/IZ
 Sample : SSTD02042
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Apr 21 19:13:04 2015
 Quant Method : Z:\HPCHEM1\BNA M\Methods\SOM01.2-EPA-BM041515.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 16 05:47:45 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
45	2,6-Dinitrotoluene	0.278	0.302	-8.6	88	0.00
46 S	Acenaphthylene-d8	1.907	1.964	-3.0	83	-0.01
47	Acenaphthylene	2.069	2.086	-0.8	81	-0.01
48	3-Nitroaniline	0.265	0.342	-29.1#	114	0.00
49 C	Acenaphthene	1.365	1.351	1.0	82	-0.01
50	2,4-Dinitrophenol	0.152	0.115	24.3#	69	0.00
51 S	4-Nitrophenol-d4	0.259	0.282	-8.9	97	0.00
52	4-Nitrophenol	0.196	0.206	-5.1	92	0.00
53	Dibenzofuran	1.898	1.912	-0.7	84	-0.01
54	2,4-Dinitrotoluene	0.412	0.442	-7.3	88	-0.01
55	2,3,4,6-Tetrachlorophenol	0.329	0.349	-6.1	89	-0.01
56	Diethylphthalate	1.480	1.509	-2.0	86	-0.01
57 S	Fluorene-d10	1.325	1.320	0.4	84	0.00
58	Fluorene	1.562	1.574	-0.8	84	-0.01
59	4-Chlorophenyl-phenylether	0.745	0.737	1.1	83	-0.01
60	4-Nitroaniline	0.329	0.359	-9.1	98	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	87	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.109	0.096	11.9	80	0.00
63	4,6-Dinitro-2-methylphenol	0.117	0.105	10.3	81	0.00
64	N-Nitrosodiphenylamine	0.616	0.620	-0.6	85	0.00
65	4-Bromophenyl-phenylether	0.194	0.198	-2.1	88	0.00
66	Hexachlorobenzene	0.219	0.219	0.0	86	0.00
67	Atrazine	0.197	0.206	-4.6	90	0.00
68 C	Pentachlorophenol	0.116	0.121	-4.3	96	0.00
69	Phenanthrene	1.120	1.132	-1.1	89	0.00
70 S	Anthracene-d10	0.905	0.938	-3.6	90	-0.01
71	Anthracene	1.140	1.164	-2.1	89	-0.01
72	Carbazole	1.013	1.068	-5.4	93	0.00
73	Di-n-butylphthalate	1.124	1.220	-8.5	95	-0.01
74 C	Fluoranthene	1.207	1.253	-3.8	93	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	96	0.00
76 S	Pyrene-d10	0.928	0.933	-0.5	93	0.00
77	Pyrene	1.270	1.275	-0.4	92	0.00
78	Butylbenzylphthalate	0.467	0.529	-13.3	108	0.00
79	3,3'-Dichlorobenzidine	0.304	0.381	-25.3#	117	0.00
80	Benzo(a)anthracene	1.163	1.207	-3.8	98	0.00
81	Bis(2-ethylhexyl)phthalate	0.698	0.793	-13.6	108	-0.01
82	Chrysene	1.108	1.127	-1.7	98	0.00
83 I	Perylene-d12	1.000	1.000	0.0	101	-0.01
84	Di-n-octyl phthalate	1.313	1.413	-7.6	109	0.00
85	Benzo(b)fluoranthene	1.234	1.173	4.9	95	0.00
86	Benzo(k)fluoranthene	1.154	1.194	-3.5	104	0.00
87 S	Benzo(a)pyrene-d12	0.895	0.906	-1.2	102	0.00

Data Path : Z:\HPCHEM1\BNA M\DATA\BM042015\
Data File : BM001099.D
Acq On : 21 Apr 2015 11:45
Operator : TP/IZ
Sample : SSTD02042
Misc :
ALS Vial : 2 Sample Multiplier: 1

Quant Time: Apr 21 19:13:04 2015
Quant Method : Z:\HPCHEM1\BNA M\Methods\SOM01.2-EPA-BM041515.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 16 05:47:45 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
88 C	Benzo(a)pyrene	1.155	1.170	-1.3	101	-0.01
89	Indeno(1,2,3-cd)pyrene	1.256	1.321	-5.2	105	0.00
90	Dibenzo(a,h)anthracene	1.057	1.106	-4.6	105	-0.02
91	Benzo(a,h,i)perylene	1.061	1.108	-4.4	104	-0.02

(#) = Out of Range

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