

Data Path : Z:\HPCHEM1\BNA M\Data\BM041516\
 Data File : BM004954.D
 Acq On : 15 Apr 2016 10:37
 Operator : UM/SJ
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02050

Quant Time: Apr 15 17:32:05 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM041416.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Apr 15 00:36:20 2016
 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	77	0.00
2	1,4-Dioxane	0.681	0.705	-3.5	78	0.00
3 S	1,4-Dioxane-d8	0.373	0.388	-4.0	78	0.00
4	Benzaldehyde	0.816	0.959	-17.5	74	0.00
5 S	Phenol-d5	1.428	1.452	-1.7	73	0.00
6	Phenol	1.478	1.508	-2.0	73	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	0.819	0.848	-3.5	75	0.00
8	Bis(2-Chloroethyl)ether	1.155	1.193	-3.3	74	0.00
9 S	2-Chlorophenol-d4	1.193	1.193	0.0	76	0.00
10	2-Chlorophenol	1.226	1.194	2.6	73	0.00
11	2-Methylphenol	1.161	1.156	0.4	72	0.00
12	2,2'-oxybis(1-Chloropropane	1.434	1.475	-2.9	73	0.00
13 S	4-Methylphenol-d8	1.179	1.172	0.6	72	0.00
14	Acetophenone	1.765	1.837	-4.1	71	0.00
15 P	N-Nitroso-di-n-propylamine	0.931	0.876	5.9	68	0.00
16	4-Methylphenol	1.257	1.251	0.5	71	0.00
17	Hexachloroethane	0.481	0.478	0.6	77	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	74	0.00
19 S	Nitrobenzene-d5	0.132	0.130	1.5	71	0.00
20	Nitrobenzene	0.309	0.311	-0.6	73	0.00
21	Isophorone	0.587	0.554	5.6	68	0.00
22 S	2-Nitrophenol-d4	0.149	0.142	4.7	69	0.00
23 C	2-Nitrophenol	0.158	0.155	1.9	69	0.00
24	2,4-Dimethylphenol	0.318	0.321	-0.9	71	0.00
25	Bis(2-Chloroethoxy)methane	0.369	0.357	3.3	70	0.00
26 S	2,4-Dichlorophenol-d3	0.271	0.264	2.6	70	0.00
27 C	2,4-Dichlorophenol	0.273	0.272	0.4	72	0.00
28	Naphthalene	0.891	0.916	-2.8	74	0.00
29 S	4-Chloroaniline-d4	0.305	0.344	-12.8	70	0.00
30	4-Chloroaniline	0.308	0.352	-14.3	71	0.00
31 C	Hexachlorobutadiene	0.176	0.181	-2.8	75	0.00
32	Caprolactam	0.084	0.073	13.1	64	0.00
33 C	4-Chloro-3-methylphenol	0.301	0.285	5.3	68	0.00
34	2-Methylnaphthalene	0.649	0.655	-0.9	71	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	70	0.00
36	1,2,4,5-Tetrachlorobenzene	0.571	0.597	-4.6	71	0.00
37	Hexachlorocyclopentadiene	0.306	0.322	-5.2	78	0.00
38 C	2,4,6-Trichlorophenol	0.376	0.368	2.1	67	0.00
39	2,4,5-Trichlorophenol	0.406	0.404	0.5	68	0.00
40	1,1'-Biphenyl	1.435	1.500	-4.5	71	0.00
41	2-Chloronaphthalene	1.091	1.137	-4.2	71	0.00
42	2-Nitroaniline	0.307	0.285	7.2	64	0.00
43 S	Dimethylphthalate-d6	1.360	1.327	2.4	67	0.00
44	Dimethylphthalate	1.354	1.323	2.3	67	0.00

Data Path : Z:\HPCHEM1\BNA M\Data\BM041516\
 Data File : BM004954.D
 Acq On : 15 Apr 2016 10:37
 Operator : UM/SJ
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02050

Quant Time: Apr 15 17:32:05 2016
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM041416.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Apr 15 00:36:20 2016
 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	0.270	0.255	5.6	64	0.00
46 S	Acenaphthylene-d8	1.668	1.697	-1.7	69	0.00
47	Acenaphthylene	1.772	1.831	-3.3	70	0.00
48	3-Nitroaniline	0.273	0.278	-1.8	67	0.00
49 C	Acenaphthene	1.153	1.178	-2.2	70	0.00
50	2,4-Dinitrophenol	0.119	0.093	21.8#	74	0.00
51 S	4-Nitrophenol-d4	0.232	0.214	7.8	67	0.00
52	4-Nitrophenol	0.196	0.182	7.1	67	0.00
53	Dibenzofuran	1.669	1.696	-1.6	69	0.00
54	2,4-Dinitrotoluene	0.384	0.370	3.6	66	0.00
55	2,3,4,6-Tetrachlorophenol	0.334	0.327	2.1	68	0.00
56	Diethylphthalate	1.326	1.280	3.5	67	0.00
57 S	Fluorene-d10	1.175	1.187	-1.0	70	0.00
58	Fluorene	1.284	1.321	-2.9	70	0.00
59	4-Chlorophenyl-phenylether	0.643	0.654	-1.7	68	0.00
60	4-Nitroaniline	0.283	0.284	-0.4	72	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	70	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.087	0.085	2.3	75	0.00
63	4,6-Dinitro-2-methylphenol	0.090	0.092	-2.2	76	0.00
64	N-Nitrosodiphenylamine	0.516	0.529	-2.5	69	0.00
65	4-Bromophenyl-phenylether	0.184	0.184	0.0	68	0.00
66	Hexachlorobenzene	0.206	0.209	-1.5	69	0.00
67	Atrazine	0.175	0.175	0.0	66	0.00
68 C	Pentachlorophenol	0.113	0.106	6.2	68	0.00
69	Phenanthrene	0.933	0.971	-4.1	71	0.00
70 S	Anthracene-d10	0.795	0.817	-2.8	70	0.00
71	Anthracene	0.944	0.987	-4.6	71	0.00
72	Carbazole	0.795	0.872	-9.7	72	0.00
73	Di-n-butylphthalate	0.957	0.919	4.0	67	0.00
74 C	Fluoranthene	0.961	1.065	-10.8	73	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	79	0.00
76 S	Pyrene-d10	0.833	0.803	3.6	74	0.00
77	Pyrene	1.057	1.041	1.5	75	0.00
78	Butylbenzylphthalate	0.409	0.369	9.8	70	0.00
79	3,3'-Dichlorobenzidine	0.293	0.335	-14.3	79	0.00
80	Benzo(a)anthracene	0.985	0.993	-0.8	79	0.00
81	Bis(2-ethylhexyl)phthalate	0.559	0.532	4.8	74	0.00
82	Chrysene	0.934	0.962	-3.0	81	0.00
83 I	Perylene-d12	1.000	1.000	0.0	81	0.00
84	Di-n-octyl phthalate	0.960	0.952	0.8	76	0.00
85	Benzo(b)fluoranthene	0.992	1.050	-5.8	84	0.00
86	Benzo(k)fluoranthene	0.957	0.947	1.0	77	0.00
87 S	Benzo(a)pyrene-d12	0.770	0.786	-2.1	81	0.00

Data Path : Z:\HPCHEM1\BNA M\Data\BM041516\
Data File : BM004954.D
Acq On : 15 Apr 2016 10:37
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02050

Quant Time: Apr 15 17:32:05 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM041416.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Apr 15 00:36:20 2016
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)
88 C	Benzo(a)pyrene	0.950	0.983	-3.5	81	0.00
89	Indeno(1,2,3-cd)pyrene	1.057	1.105	-4.5	82	0.00
90	Dibenzo(a,h)anthracene	0.885	0.934	-5.5	82	0.00
91	Benzo(a,h,i)perylene	0.890	0.929	-4.4	82	0.00

(#) = Out of Range

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