

Data Path : Z:\HPCHEM1\BNA_M\Data\BM012015\
 Data File : BM000304.D
 Acq On : 21 Jan 2015 17:02
 Operator : TP/IZ
 Sample : SSTD02041
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 LabSampled :
 SSTD02041

Quant Time: Jan 22 04:10:51 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM122914.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 22 04:03:17 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)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	57	-0.04
2	1,4-Dioxane	0.478	0.607	-27.0#	75	-0.02
3 S	1,4-Dioxane-d8	0.345	0.434	-25.8#	73	-0.02
4	Benzaldehyde	1.172	1.335	-13.9	62	-0.02
5 S	Phenol-d5	1.682	1.551	7.8	54	-0.03
6	Phenol	1.774	1.679	5.4	55	-0.02
7 S	Bis-(2-Chloroethyl)ether-d8	1.063	1.119	-5.3	63	-0.03
8	Bis(2-Chloroethyl)ether	1.377	1.372	0.4	59	-0.02
9 S	2-Chlorophenol-d4	1.234	1.186	3.9	57	-0.02
10	2-Chlorophenol	1.291	1.188	8.0	55	-0.03
11	2-Methylphenol	1.366	1.217	10.9	52	-0.02
12	2,2'-oxybis(1-Chloropropane	2.264	2.447	-8.1	64	-0.03
13 S	4-Methylphenol-d8	1.338	1.176	12.1	52	-0.02
14	Acetophenone	2.352	2.233	5.1	56	-0.03
15 P	N-Nitroso-di-n-propylamine	1.223	1.133	7.4	54	-0.03
16	4-Methylphenol	1.475	1.292	12.4	51	-0.03
17	Hexachloroethane	0.590	0.654	-10.8	67	-0.04
18 I	Naphthalene-d8	1.000	1.000	0.0	52	-0.03
19 S	Nitrobenzene-d5	0.138	0.151	-9.4	59	-0.03
20	Nitrobenzene	0.412	0.501	-21.6#	64	-0.03
21	Isophorone	0.795	0.783	1.5	52	-0.03
22 S	2-Nitrophenol-d4	0.141	0.148	-5.0	56	-0.04
23 C	2-Nitrophenol	0.156	0.158	-1.3	52	-0.03
24	2,4-Dimethylphenol	0.423	0.412	2.6	51	-0.03
25	Bis(2-Chloroethoxy)methane	0.442	0.429	2.9	52	-0.03
26 S	2,4-Dichlorophenol-d3	0.312	0.296	5.1	51	-0.03
27 C	2,4-Dichlorophenol	0.311	0.293	5.8	49	-0.04
28	Naphthalene	1.020	1.007	1.3	53	-0.03
29 S	4-Chloroaniline-d4	0.354	0.366	-3.4	48	-0.03
30	4-Chloroaniline	0.364	0.378	-3.8	48	-0.03
31 C	Hexachlorobutadiene	0.221	0.244	-10.4	59	-0.03
32	Caprolactam	0.101	0.075	25.7#	38	-0.03
33 C	4-Chloro-3-methylphenol	0.382	0.339	11.3	47	-0.02
34	2-Methylnaphthalene	0.756	0.713	5.7	50	-0.03
35 I	Acenaphthene-d10	1.000	1.000	0.0	47	-0.03
36	1,2,4,5-Tetrachlorobenzene	0.555	0.597	-7.6	53	-0.03
37	Hexachlorocyclopentadiene	0.345	0.373	-8.1	53	-0.03
38 C	2,4,6-Trichlorophenol	0.354	0.339	4.2	47	-0.02
39	2,4,5-Trichlorophenol	0.384	0.363	5.5	45	-0.02
40	1,1'-Biphenyl	1.441	1.444	-0.2	48	-0.03
41	2-Chloronaphthalene	1.105	1.119	-1.3	50	-0.02
42	2-Nitroaniline	0.377	0.359	4.8	48	-0.02
43 S	Dimethylphthalate-d6	1.475	1.380	6.4	45	-0.03
44	Dimethylphthalate	1.469	1.370	6.7	45	-0.02

Data Path : Z:\HPCHEM1\BNA_M\Data\BM012015\
 Data File : BM000304.D
 Acq On : 21 Jan 2015 17:02
 Operator : TP/IZ
 Sample : SSTD02041
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 LabSampled :
 SSTD02041

Quant Time: Jan 22 04:10:51 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM122914.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 22 04:03:17 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)
45	2,6-Dinitrotoluene	0.263	0.255	3.0	46	-0.03
46 S	Acenaphthylene-d8	1.767	1.739	1.6	48	-0.03
47	Acenaphthylene	1.863	1.829	1.8	47	-0.03
48	3-Nitroaniline	0.279	0.256	8.2	43	-0.02
49 C	Acenaphthene	1.188	1.174	1.2	48	-0.04
50	2,4-Dinitrophenol	0.128	0.095	25.8#	40	-0.02
51 S	4-Nitrophenol-d4	0.240	0.194	19.2	40	-0.02
52	4-Nitrophenol	0.318	0.307	3.5	48	-0.01
53	Dibenzofuran	1.720	1.723	-0.2	49	-0.03
54	2,4-Dinitrotoluene	0.392	0.396	-1.0	49	-0.02
55	2,3,4,6-Tetrachlorophenol	0.331	0.312	5.7	44	-0.02
56	Diethylphthalate	1.544	1.438	6.9	45	-0.03
57 S	Fluorene-d10	1.267	1.206	4.8	46	-0.03
58	Fluorene	1.427	1.392	2.5	47	-0.03
59	4-Chlorophenyl-phenylether	0.704	0.705	-0.1	49	-0.02
60	4-Nitroaniline	0.312	0.276	11.5	43	-0.02
61 I	Phenanthrene-d10	1.000	1.000	0.0	44	-0.03
62 S	4,6-Dinitro-2-methylphenol-	0.095	0.082	13.7	43	-0.03
63	4,6-Dinitro-2-methylphenol	0.098	0.089	9.2	44	-0.02
64	N-Nitrosodiphenylamine	0.580	0.579	0.2	45	-0.03
65	4-Bromophenyl-phenylether	0.204	0.214	-4.9	48	-0.03
66	Hexachlorobenzene	0.235	0.250	-6.4	49	-0.02
67	Atrazine	0.226	0.197	12.8	39	-0.03
68 C	Pentachlorophenol	0.142	0.118	16.9	39	-0.02
69	Phenanthrene	1.083	1.135	-4.8	47	-0.03
70 S	Anthracene-d10	0.925	0.952	-2.9	46	-0.02
71	Anthracene	1.113	1.148	-3.1	46	-0.02
72	Carbazole	1.006	0.990	1.6	44	-0.02
73	Di-n-butylphthalate	1.196	1.141	4.6	41	-0.03
74 C	Fluoranthene	1.265	1.304	-3.1	45	-0.03
75 I	Chrysene-d12	1.000	1.000	0.0	45	-0.02
76 S	Pyrene-d10	0.923	0.886	4.0	45	-0.02
77	Pyrene	1.205	1.192	1.1	46	-0.03
78	Butylbenzylphthalate	0.490	0.390	20.4#	36	-0.03
79	3,3'-Dichlorobenzidine	0.353	0.301	14.7	35	-0.03
80	Benzo(a)anthracene	1.155	1.157	-0.2	47	-0.02
81	Bis(2-ethylhexyl)phthalate	0.713	0.603	15.4	38	-0.03
82	Chrysene	1.064	1.111	-4.4	47	-0.02
83 I	Perylene-d12	1.000	1.000	0.0	48	-0.03
84	Di-n-octyl phthalate	1.298	1.099	15.3	40	-0.05
85	Benzo(b)fluoranthene	1.276	1.180	7.5	44	-0.03
86	Benzo(k)fluoranthene	1.090	1.180	-8.3	54	-0.04
87 S	Benzo(a)pyrene-d12	0.908	0.921	-1.4	50	-0.04

Data Path : Z:\HPCHEM1\BNA_M\Data\BM012015\
Data File : BM000304.D
Acq On : 21 Jan 2015 17:02
Operator : TP/IZ
Sample : SSTD02041
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
LabSampleId :
SSTD02041

Quant Time: Jan 22 04:10:51 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-BM122914.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jan 22 04:03:17 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	1.121	1.163	-3.7	50	-0.04
89	Indeno(1,2,3-cd)pyrene	1.170	1.342	-14.7	58	-0.06
90	Dibenzo(a,h)anthracene	0.985	1.124	-14.1	57	-0.06
91	Benzo(g,h,i)perylene	0.961	1.130	-17.6	60	-0.07

(#) = Out of Range

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