

Data Path : Z:\HPCHEM1\BNA M\Data\BM022118\
 Data File : BM014318.D
 Acq On : 21 Feb 2018 09:23
 Operator : SJ/JU
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02048

Quant Time: Feb 22 03:32:12 2018
 Quant Method : Z:\HPCHEM1\BNA M\Methods\SOM-EPA-BM021418.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Feb 20 03:49:19 2018
 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	82	0.00
2	1,4-Dioxane	0.519	0.579	-11.6	95	0.00
3 S	1,4-Dioxane-d8	0.467	0.488	-4.5	91	0.00
4	Benzaldehyde	0.689	0.780	-13.2	80	0.00
5 S	Phenol-d5	1.690	1.656	2.0	86	0.00
6	Phenol	1.774	1.718	3.2	86	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	1.023	1.074	-5.0	88	0.00
8	Bis(2-Chloroethyl)ether	1.287	1.215	5.6	80	0.00
9 S	2-Chlorophenol-d4	1.296	1.357	-4.7	86	0.00
10	2-Chlorophenol	1.323	1.289	2.6	86	0.00
11	2-Methylphenol	1.357	1.332	1.8	88	0.00
12	2,2'-oxybis(1-Chloropropane	1.299	1.224	5.8	83	0.00
13 S	4-Methylphenol-d8	1.479	1.425	3.7	85	0.00
14	Acetophenone	2.552	2.524	1.1	90	0.00
15 P	N-Nitroso-di-n-propylamine	1.279	1.372	-7.3	94	0.00
16	4-Methylphenol	1.478	1.400	5.3	84	0.00
17	Hexachloroethane	0.689	0.720	-4.5	88	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	86	0.00
19 S	Nitrobenzene-d5	0.148	0.140	5.4	83	0.00
20	Nitrobenzene	0.437	0.435	0.5	89	0.00
21	Isophorone	0.748	0.736	1.6	88	0.00
22 S	2-Nitrophenol-d4	0.153	0.149	2.6	82	0.00
23 C	2-Nitrophenol	0.166	0.169	-1.8	88	0.00
24	2,4-Dimethylphenol	0.406	0.408	-0.5	91	0.00
25	Bis(2-Chloroethoxy)methane	0.394	0.392	0.5	87	0.00
26 S	2,4-Dichlorophenol-d3	0.317	0.300	5.4	82	0.00
27 C	2,4-Dichlorophenol	0.301	0.283	6.0	86	0.00
28	Naphthalene	1.027	1.011	1.6	88	0.00
29 S	4-Chloroaniline-d4	0.312	0.348	-11.5	92	0.00
30	4-Chloroaniline	0.318	0.350	-10.1	90	0.00
31 C	Hexachlorobutadiene	0.239	0.249	-4.2	95	0.00
32	Caprolactam	0.096	0.094	2.1	87	0.00
33 C	4-Chloro-3-methylphenol	0.363	0.364	-0.3	89	0.00
34	2-Methylnaphthalene	0.782	0.804	-2.8	91	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	91	0.00
36	1,2,4,5-Tetrachlorobenzene	0.642	0.623	3.0	91	0.00
37	Hexachlorocyclopentadiene	0.349	0.302	13.5	92	0.00
38 C	2,4,6-Trichlorophenol	0.401	0.397	1.0	93	0.00
39	2,4,5-Trichlorophenol	0.409	0.370	9.5	78	0.00
40	1,1'-Biphenyl	1.541	1.528	0.8	90	0.00
41	2-Chloronaphthalene	1.231	1.213	1.5	90	0.00
42	2-Nitroaniline	0.416	0.419	-0.7	91	0.00
43 S	Dimethylphthalate-d6	1.652	1.637	0.9	92	0.00
44	Dimethylphthalate	1.629	1.597	2.0	92	0.00

Data Path : Z:\HPCHEM1\BNA M\Data\BM022118\
 Data File : BM014318.D
 Acq On : 21 Feb 2018 09:23
 Operator : SJ/JU
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02048

Quant Time: Feb 22 03:32:12 2018
 Quant Method : Z:\HPCHEM1\BNA M\Methods\SOM-EPA-BM021418.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Feb 20 03:49:19 2018
 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	0.309	0.316	-2.3	94	0.00
46 S	Acenaphthylene-d8	2.037	2.057	-1.0	94	0.00
47	Acenaphthylene	1.925	1.932	-0.4	92	0.00
48	3-Nitroaniline	0.262	0.232	11.5	92	0.00
49 C	Acenaphthene	1.355	1.366	-0.8	95	0.00
50	2,4-Dinitrophenol	0.166	0.119	28.3#	90	0.00
51 S	4-Nitrophenol-d4	0.228	0.176	22.8#	88	0.00
52	4-Nitrophenol	0.299	0.229	23.4#	80	0.00
53	Dibenzofuran	2.032	2.089	-2.8	95	0.00
54	2,4-Dinitrotoluene	0.496	0.520	-4.8	96	0.00
55	2,3,4,6-Tetrachlorophenol	0.414	0.416	-0.5	93	0.00
56	Diethylphthalate	1.815	1.826	-0.6	93	0.00
57 S	Fluorene-d10	1.480	1.514	-2.3	95	0.00
58	Fluorene	1.755	1.761	-0.3	94	0.00
59	4-Chlorophenyl-phenylether	0.918	0.927	-1.0	95	0.00
60	4-Nitroaniline	0.294	0.238	19.0	90	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	91	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.120	0.105	12.5	93	0.00
63	4,6-Dinitro-2-methylphenol	0.122	0.104	14.8	88	0.00
64	N-Nitrosodiphenylamine	0.575	0.562	2.3	93	0.00
65	4-Bromophenyl-phenylether	0.223	0.224	-0.4	97	0.00
66	Hexachlorobenzene	0.235	0.237	-0.9	99	0.00
67	Atrazine	0.227	0.226	0.4	97	0.00
68 C	Pentachlorophenol	0.123	0.111	9.8	99	0.00
69	Phenanthrene	1.152	1.142	0.9	95	0.00
70 S	Anthracene-d10	0.965	0.974	-0.9	95	0.00
71	Anthracene	1.159	1.151	0.7	92	0.00
72	Carbazole	0.974	0.909	6.7	92	0.00
73	Di-n-butylphthalate	1.257	1.262	-0.4	93	0.00
74 C	Fluoranthene	1.377	1.372	0.4	96	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	92	0.00
76 S	Pyrene-d10	1.032	1.022	1.0	95	0.00
77	Pyrene	1.329	1.293	2.7	93	0.00
78	Butylbenzylphthalate	0.550	0.539	2.0	92	0.00
79	3,3'-Dichlorobenzidine	0.352	0.317	9.9	99	0.00
80	Benzo(a)anthracene	1.248	1.252	-0.3	96	0.00
81	Bis(2-ethylhexyl)phthalate	0.787	0.776	1.4	93	0.00
82	Chrysene	1.164	1.177	-1.1	96	0.00
83 I	Perylene-d12	1.000	1.000	0.0	99	0.00
84	Di-n-octyl phthalate	1.716	1.486	13.4	96	0.00
85	Benzo(b)fluoranthene	1.339	1.352	-1.0	101	0.00
86	Benzo(k)fluoranthene	1.273	1.249	1.9	100	0.00
87 S	Benzo(a)pyrene-d12	0.973	0.969	0.4	100	0.00

Data Path : Z:\HPCHEM1\BNA M\Data\BM022118\
Data File : BM014318.D
Acq On : 21 Feb 2018 09:23
Operator : SJ/JU
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02048

Quant Time: Feb 22 03:32:12 2018
Quant Method : Z:\HPCHEM1\BNA M\Methods\SOM-EPA-BM021418.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Feb 20 03:49:19 2018
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	1.197	1.167	2.5	101	0.00
89	Indeno(1,2,3-cd)pyrene	1.161	1.123	3.3	101	0.00
90	Dibenzo(a,h)anthracene	0.965	0.931	3.5	101	0.00
91	Benzo(a,h,i)perylene	0.941	0.909	3.4	103	0.00

(#) = Out of Range

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