

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM082916\
 Data File : BM007326.D
 Acq On : 30 Aug 2016 01:34
 Operator : UM/SJ
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02036

Quant Time: Aug 30 04:01:40 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM082316.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 30 02:26:11 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	69	0.00
2	1,4-Dioxane	0.436	0.417	4.4	67	0.00
3 S	1,4-Dioxane-d8	0.391	0.443	-13.3	79	0.00
4	Benzaldehyde	1.076	1.189	-10.5	67	0.00
5 S	Phenol-d5	1.889	1.838	2.7	67	0.00
6	Phenol	1.960	1.955	0.3	68	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	0.993	1.028	-3.5	69	0.00
8	Bis(2-Chloroethyl)ether	1.365	1.389	-1.8	68	0.00
9 S	2-Chlorophenol-d4	1.396	1.383	0.9	68	0.00
10	2-Chlorophenol	1.430	1.411	1.3	66	0.00
11	2-Methylphenol	1.532	1.509	1.5	68	0.00
12	2,2'-oxybis(1-Chloropropane	1.536	1.564	-1.8	67	0.00
13 S	4-Methylphenol-d8	1.648	1.643	0.3	68	0.00
14	Acetophenone	2.527	2.624	-3.8	69	0.00
15 P	N-Nitroso-di-n-propylamine	1.343	1.385	-3.1	68	0.00
16	4-Methylphenol	1.718	1.743	-1.5	69	0.00
17	Hexachloroethane	0.556	0.581	-4.5	71	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	69	0.00
19 S	Nitrobenzene-d5	0.147	0.150	-2.0	69	0.00
20	Nitrobenzene	0.377	0.393	-4.2	70	0.00
21	Isophorone	0.768	0.779	-1.4	67	0.00
22 S	2-Nitrophenol-d4	0.171	0.170	0.6	67	0.00
23 C	2-Nitrophenol	0.186	0.191	-2.7	67	0.00
24	2,4-Dimethylphenol	0.411	0.435	-5.8	71	0.00
25	Bis(2-Chloroethoxy)methane	0.433	0.439	-1.4	67	0.00
26 S	2,4-Dichlorophenol-d3	0.341	0.351	-2.9	69	0.00
27 C	2,4-Dichlorophenol	0.339	0.350	-3.2	68	0.00
28	Naphthalene	0.995	1.013	-1.8	68	0.00
29 S	4-Chloroaniline-d4	0.426	0.460	-8.0	69	0.00
30	4-Chloroaniline	0.438	0.474	-8.2	68	0.00
31 C	Hexachlorobutadiene	0.229	0.245	-7.0	71	0.00
32	Caprolactam	0.137	0.140	-2.2	66	0.00
33 C	4-Chloro-3-methylphenol	0.420	0.443	-5.5	69	0.00
34	2-Methylnaphthalene	0.814	0.854	-4.9	70	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	70	0.00
36	1,2,4,5-Tetrachlorobenzene	0.666	0.686	-3.0	70	0.00
37	Hexachlorocyclopentadiene	0.402	0.386	4.0	68	0.00
38 C	2,4,6-Trichlorophenol	0.471	0.470	0.2	68	0.00
39	2,4,5-Trichlorophenol	0.491	0.501	-2.0	70	0.00
40	1,1'-Biphenyl	1.551	1.600	-3.2	71	0.00
41	2-Chloronaphthalene	1.216	1.232	-1.3	69	0.00
42	2-Nitroaniline	0.397	0.420	-5.8	71	0.00
43 S	Dimethylphthalate-d6	1.725	1.789	-3.7	71	0.00
44	Dimethylphthalate	1.719	1.798	-4.6	71	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM082916\
 Data File : BM007326.D
 Acq On : 30 Aug 2016 01:34
 Operator : UM/SJ
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02036

Quant Time: Aug 30 04:01:40 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM082316.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 30 02:26:11 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	2,6-Dinitrotoluene	0.351	0.365	-4.0	70	0.00
46 S	Acenaphthylene-d8	2.000	2.077	-3.8	71	0.00
47	Acenaphthylene	2.058	2.132	-3.6	70	0.00
48	3-Nitroaniline	0.359	0.382	-6.4	68	0.00
49 C	Acenaphthene	1.339	1.377	-2.8	70	0.00
50	2,4-Dinitrophenol	0.238	0.224	5.9	70	0.00
51 S	4-Nitrophenol-d4	0.326	0.338	-3.7	70	0.00
52	4-Nitrophenol	0.347	0.377	-8.6	74	0.00
53	Dibenzofuran	2.004	2.088	-4.2	71	0.00
54	2,4-Dinitrotoluene	0.538	0.573	-6.5	72	0.00
55	2,3,4,6-Tetrachlorophenol	0.504	0.524	-4.0	70	0.00
56	Diethylphthalate	1.793	1.885	-5.1	71	0.00
57 S	Fluorene-d10	1.492	1.559	-4.5	71	0.00
58	Fluorene	1.650	1.750	-6.1	72	0.00
59	4-Chlorophenyl-phenylether	0.863	0.917	-6.3	73	0.00
60	4-Nitroaniline	0.407	0.428	-5.2	68	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	74	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.126	0.122	3.2	71	0.00
63	4,6-Dinitro-2-methylphenol	0.134	0.133	0.7	72	0.00
64	N-Nitrosodiphenylamine	0.558	0.563	-0.9	71	0.00
65	4-Bromophenyl-phenylether	0.226	0.229	-1.3	71	0.00
66	Hexachlorobenzene	0.253	0.256	-1.2	72	0.00
67	Atrazine	0.236	0.245	-3.8	70	0.00
68 C	Pentachlorophenol	0.163	0.159	2.5	70	0.00
69	Phenanthrene	1.074	1.121	-4.4	74	0.00
70 S	Anthracene-d10	0.934	0.963	-3.1	73	0.00
71	Anthracene	1.108	1.152	-4.0	73	0.00
72	Carbazole	0.963	1.025	-6.4	72	0.00
73	Di-n-butylphthalate	1.212	1.218	-0.5	70	0.00
74 C	Fluoranthene	1.347	1.512	-12.2	75	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	75	0.00
76 S	Pyrene-d10	0.836	0.858	-2.6	75	0.00
77	Pyrene	1.073	1.106	-3.1	75	0.00
78	Butylbenzylphthalate	0.446	0.440	1.3	72	0.00
79	3,3'-Dichlorobenzidine	0.397	0.420	-5.8	69	0.00
80	Benzo(a)anthracene	1.179	1.235	-4.7	76	0.00
81	Bis(2-ethylhexyl)phthalate	0.648	0.644	0.6	72	0.00
82	Chrysene	1.068	1.105	-3.5	75	0.00
83 I	Perylene-d12	1.000	1.000	0.0	73	0.00
84	Di-n-octyl phthalate	0.996	1.073	-7.7	70	0.00
85	Benzo(b)fluoranthene	1.183	1.228	-3.8	71	0.00
86	Benzo(k)fluoranthene	1.093	1.197	-9.5	77	0.00
87 S	Benzo(a)pyrene-d12	0.921	0.949	-3.0	73	0.00

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM082916\
Data File : BM007326.D
Acq On : 30 Aug 2016 01:34
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02036

Quant Time: Aug 30 04:01:40 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM082316.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 30 02:26:11 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
88 C	Benzo(a)pyrene	1.132	1.190	-5.1	73	0.00
89	Indeno(1,2,3-cd)pyrene	1.389	1.382	0.5	69	0.00
90	Dibenzo(a,h)anthracene	1.155	1.155	0.0	69	0.00
91	Benzo(g,h,i)perylene	1.179	1.147	2.7	68	0.00

(#) = Out of Range

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