

Data Path : Z:\HPCHEM1\BNA_M\Data\BM081916\
 Data File : BM007207.D
 Acq On : 18 Aug 2016 17:03
 Operator : UM/SJ
 Sample : SSTDC0020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02052

Quant Time: Aug 19 04:27:33 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM081116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 19 04:26:41 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	92	0.00
2	1,4-Dioxane	0.476	0.487	-2.3	96	0.00
3 S	1,4-Dioxane-d8	0.458	0.486	-6.1	96	0.00
4	Benzaldehyde	0.974	0.994	-2.1	87	0.00
5 S	Phenol-d5	1.620	1.580	2.5	91	0.00
6	Phenol	1.707	1.660	2.8	90	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	0.929	0.906	2.5	89	0.00
8	Bis(2-Chloroethyl)ether	1.259	1.250	0.7	90	0.00
9 S	2-Chlorophenol-d4	1.290	1.296	-0.5	93	0.00
10	2-Chlorophenol	1.320	1.307	1.0	90	0.00
11	2-Methylphenol	1.283	1.211	5.6	89	0.00
12	2,2'-oxybis(1-Chloropropane	1.443	1.390	3.7	90	0.00
13 S	4-Methylphenol-d8	1.333	1.239	7.1	88	0.00
14	Acetophenone	2.053	1.949	5.1	87	0.00
15 P	N-Nitroso-di-n-propylamine	1.068	0.980	8.2	85	0.00
16	4-Methylphenol	1.401	1.316	6.1	89	0.00
17	Hexachloroethane	0.551	0.546	0.9	91	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	90	0.00
19 S	Nitrobenzene-d5	0.145	0.148	-2.1	88	0.00
20	Nitrobenzene	0.374	0.386	-3.2	91	0.00
21	Isophorone	0.700	0.670	4.3	84	0.00
22 S	2-Nitrophenol-d4	0.155	0.164	-5.8	92	0.00
23 C	2-Nitrophenol	0.170	0.179	-5.3	91	0.00
24	2,4-Dimethylphenol	0.389	0.392	-0.8	88	0.00
25	Bis(2-Chloroethoxy)methane	0.420	0.412	1.9	86	0.00
26 S	2,4-Dichlorophenol-d3	0.325	0.331	-1.8	90	0.00
27 C	2,4-Dichlorophenol	0.324	0.327	-0.9	89	0.00
28	Naphthalene	0.991	1.003	-1.2	89	0.00
29 S	4-Chloroaniline-d4	0.387	0.386	0.3	84	0.00
30	4-Chloroaniline	0.397	0.397	0.0	84	0.00
31 C	Hexachlorobutadiene	0.239	0.248	-3.8	91	0.00
32	Caprolactam	0.110	0.091	17.3	76	0.00
33 C	4-Chloro-3-methylphenol	0.354	0.340	4.0	84	0.00
34	2-Methylnaphthalene	0.764	0.748	2.1	86	0.00
35 I	Acenaphthene-d10	1.000	1.000	0.0	83	0.00
36	1,2,4,5-Tetrachlorobenzene	0.707	0.770	-8.9	86	0.00
37	Hexachlorocyclopentadiene	0.450	0.342	24.0#	61	0.00
38 C	2,4,6-Trichlorophenol	0.454	0.471	-3.7	82	0.00
39	2,4,5-Trichlorophenol	0.475	0.496	-4.4	82	0.00
40	1,1'-Biphenyl	1.605	1.681	-4.7	84	0.00
41	2-Chloronaphthalene	1.259	1.302	-3.4	83	0.00
42	2-Nitroaniline	0.361	0.365	-1.1	80	0.00
43 S	Dimethylphthalate-d6	1.653	1.580	4.4	77	0.00
44	Dimethylphthalate	1.652	1.587	3.9	78	0.00

Data Path : Z:\HPCHEM1\BNA_M\Data\BM081916\
 Data File : BM007207.D
 Acq On : 18 Aug 2016 17:03
 Operator : UM/SJ
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02052

Quant Time: Aug 19 04:27:33 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM081116.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Aug 19 04:26:41 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	0.321	0.325	-1.2	81	0.00
46 S	Acenaphthylene-d8	1.985	2.004	-1.0	80	0.00
47	Acenaphthylene	2.043	2.084	-2.0	81	0.00
48	3-Nitroaniline	0.317	0.317	0.0	76	0.00
49 C	Acenaphthene	1.327	1.338	-0.8	81	0.00
50	2,4-Dinitrophenol	0.190	0.130	31.6#	64	0.00
51 S	4-Nitrophenol-d4	0.296	0.259	12.5	72	0.00
52	4-Nitrophenol	0.301	0.275	8.6	75	0.00
53	Dibenzofuran	1.969	1.975	-0.3	81	0.00
54	2,4-Dinitrotoluene	0.474	0.460	3.0	78	0.00
55	2,3,4,6-Tetrachlorophenol	0.459	0.457	0.4	79	0.00
56	Diethylphthalate	1.742	1.590	8.7	75	0.00
57 S	Fluorene-d10	1.445	1.425	1.4	80	0.00
58	Fluorene	1.588	1.571	1.1	79	0.00
59	4-Chlorophenyl-phenylether	0.838	0.823	1.8	79	0.00
60	4-Nitroaniline	0.372	0.335	9.9	71	0.00
61 I	Phenanthrene-d10	1.000	1.000	0.0	78	0.00
62 S	4,6-Dinitro-2-methylphenol-	0.111	0.094	15.3	69	0.00
63	4,6-Dinitro-2-methylphenol	0.120	0.099	17.5	66	0.00
64	N-Nitrosodiphenylamine	0.566	0.582	-2.8	76	0.00
65	4-Bromophenyl-phenylether	0.227	0.233	-2.6	79	0.00
66	Hexachlorobenzene	0.256	0.262	-2.3	78	0.00
67	Atrazine	0.228	0.225	1.3	73	0.00
68 C	Pentachlorophenol	0.158	0.150	5.1	75	0.00
69	Phenanthrene	1.076	1.083	-0.7	76	0.00
70 S	Anthracene-d10	0.927	0.950	-2.5	77	0.00
71	Anthracene	1.104	1.123	-1.7	76	0.00
72	Carbazole	0.951	0.982	-3.3	75	0.00
73	Di-n-butylphthalate	1.163	1.152	0.9	74	0.00
74 C	Fluoranthene	1.305	1.390	-6.5	76	0.00
75 I	Chrysene-d12	1.000	1.000	0.0	75	0.00
76 S	Pyrene-d10	0.844	0.856	-1.4	76	0.00
77	Pyrene	1.106	1.111	-0.5	75	0.00
78	Butylbenzylphthalate	0.430	0.427	0.7	73	0.00
79	3,3'-Dichlorobenzidine	0.386	0.386	0.0	64	0.00
80	Benzo(a)anthracene	1.182	1.204	-1.9	75	0.00
81	Bis(2-ethylhexyl)phthalate	0.636	0.597	6.1	68	0.00
82	Chrysene	1.080	1.075	0.5	73	0.00
83 I	Perylene-d12	1.000	1.000	0.0	73	0.00
84	Di-n-octyl phthalate	1.032	1.026	0.6	69	0.00
85	Benzo(b)fluoranthene	1.205	1.206	-0.1	71	0.00
86	Benzo(k)fluoranthene	1.131	1.176	-4.0	75	0.00
87 S	Benzo(a)pyrene-d12	0.915	0.933	-2.0	73	0.00

Data Path : Z:\HPCHEM1\BNA_M\Data\BM081916\
Data File : BM007207.D
Acq On : 18 Aug 2016 17:03
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02052

Quant Time: Aug 19 04:27:33 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM081116.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Aug 19 04:26:41 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	1.142	1.164	-1.9	73	0.00
89	Indeno(1,2,3-cd)pyrene	1.400	1.399	0.1	71	0.01
90	Dibenzo(a,h)anthracene	1.175	1.174	0.1	71	0.01
91	Benzo(g,h,i)perylene	1.183	1.142	3.5	70	0.00

(#) = Out of Range

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