

Data Path : Z:\HPCHEM1\BNA_M\Data\BM080315\
 Data File : BM002193.D
 Acq On : 03 Aug 2015 12:24
 Operator : TP
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02094

Quant Time: Aug 04 02:31:44 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 04 02:31:26 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	80	0.00
2	1,4-Dioxane	0.413	0.374	9.4	75	0.00
3 S	1,4-Dioxane-d8	0.387	0.337	12.9	70	0.00
4	Benzaldehyde	0.759	0.766	-0.9	79	0.00
5 S	Phenol-d5	1.430	1.382	3.4	78	0.00
6	Phenol	1.462	1.418	3.0	78	0.00
7 S	Bis-(2-Chloroethyl)ether-d8	0.791	0.724	8.5	74	0.00
8	Bis(2-Chloroethyl)ether	1.099	1.026	6.6	76	0.00
9 S	2-Chlorophenol-d4	1.212	1.170	3.5	77	0.00
10	2-Chlorophenol	1.215	1.172	3.5	78	0.00
11	2-Methylphenol	1.107	1.086	1.9	80	0.00
12	2,2'-oxybis(1-Chloropropane	1.126	0.993	11.8	71	0.00
13 S	4-Methylphenol-d8	1.137	1.140	-0.3	82	0.00
14	Acetophenone	1.800	1.783	0.9	80	0.00
15 P	N-Nitroso-di-n-propylamine	0.872	0.878	-0.7	80	0.00
16	4-Methylphenol	1.189	1.197	-0.7	82	0.00
17	Hexachloroethane	0.497	0.454	8.7	75	0.00
18 I	Naphthalene-d8	1.000	1.000	0.0	84	0.00
19 S	Nitrobenzene-d5	0.142	0.135	4.9	81	0.00
20	Nitrobenzene	0.334	0.308	7.8	78	0.00
21	Isophorone	0.595	0.608	-2.2	85	0.00
22 S	2-Nitrophenol-d4	0.158	0.160	-1.3	85	0.00
23 C	2-Nitrophenol	0.173	0.166	4.0	80	0.00
24	2,4-Dimethylphenol	0.339	0.329	2.9	82	0.00
25	Bis(2-Chloroethoxy)methane	0.362	0.337	6.9	79	0.00
26 S	2,4-Dichlorophenol-d3	0.304	0.312	-2.6	85	0.00
27 C	2,4-Dichlorophenol	0.295	0.293	0.7	84	0.00
28	Naphthalene	0.950	0.884	6.9	78	0.00
29 S	4-Chloroaniline-d4	0.334	0.378	-13.2	92	0.00
30	4-Chloroaniline	0.348	0.394	-13.2	92	0.00
31 C	Hexachlorobutadiene	0.220	0.204	7.3	78	0.00
32	Caprolactam	0.083	0.102	-22.9#	107	0.00
33 C	4-Chloro-3-methylphenol	0.293	0.313	-6.8	90	0.00
34	2-Methylnaphthalene	0.691	0.686	0.7	84	0.00
35	1-Methylnaphthalene	0.665	0.655	1.5	83	0.00
36 I	Acenaphthene-d10	1.000	1.000	0.0	95	0.00
37	1,2,4,5-Tetrachlorobenzene	0.696	0.613	11.9	85	0.00
38	Hexachlorocyclopentadiene	0.396	0.269	32.1#	72	0.00
39 C	2,4,6-Trichlorophenol	0.408	0.403	1.2	93	0.00
40	2,4,5-Trichlorophenol	0.441	0.432	2.0	93	0.00
41	1,1'-Biphenyl	1.541	1.391	9.7	86	0.00
42	2-Chloronaphthalene	1.198	1.083	9.6	86	0.00
43	2-Nitroaniline	0.299	0.303	-1.3	95	0.00
44 S	Dimethylphthalate-d6	1.465	1.470	-0.3	94	0.00

Data Path : Z:\HPCHEM1\BNA_M\Data\BM080315\
 Data File : BM002193.D
 Acq On : 03 Aug 2015 12:24
 Operator : TP
 Sample : SSTDCCC020
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTD02094

Quant Time: Aug 04 02:31:44 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Aug 04 02:31:26 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	Dimethylphthalate	1.463	1.467	-0.3	94	0.00
46	2,6-Dinitrotoluene	0.298	0.310	-4.0	97	0.00
47 S	Acenaphthylene-d8	1.832	1.737	5.2	90	0.00
48	Acenaphthylene	1.890	1.768	6.5	88	0.00
49	3-Nitroaniline	0.265	0.305	-15.1	111	0.00
50 C	Acenaphthene	1.239	1.150	7.2	89	0.00
51	2,4-Dinitrophenol	0.175	0.151	13.7	94	0.00
52 S	4-Nitrophenol-d4	0.245	0.224	8.6	90	0.00
53	4-Nitrophenol	0.204	0.196	3.9	94	0.00
54	Dibenzofuran	1.777	1.697	4.5	91	0.00
55	2,4-Dinitrotoluene	0.428	0.457	-6.8	99	0.00
56	2,3,4,6-Tetrachlorophenol	0.371	0.356	4.0	89	0.00
57	Diethylphthalate	1.443	1.472	-2.0	96	0.00
58 S	Fluorene-d10	1.289	1.262	2.1	92	0.00
59	Fluorene	1.469	1.444	1.7	93	0.00
60	4-Chlorophenyl-phenylether	0.786	0.767	2.4	93	0.00
61	4-Nitroaniline	0.282	0.308	-9.2	102	0.00
62 I	Phenanthrene-d10	1.000	1.000	0.0	103	0.00
63 S	4,6-Dinitro-2-methylphenol-	0.122	0.112	8.2	103	0.00
64	4,6-Dinitro-2-methylphenol	0.128	0.116	9.4	99	0.00
65	N-Nitrosodiphenylamine	0.572	0.524	8.4	97	0.00
66	4-Bromophenyl-phenylether	0.213	0.196	8.0	96	0.00
67	Hexachlorobenzene	0.233	0.216	7.3	97	0.00
68	Atrazine	0.221	0.224	-1.4	106	0.00
69 C	Pentachlorophenol	0.119	0.082	31.1#	77	0.00
70	Phenanthrene	1.036	0.969	6.5	97	0.00
71 S	Anthracene-d10	0.893	0.843	5.6	98	0.00
72	Anthracene	1.059	0.998	5.8	99	0.00
73	1,2,3,4-Tetrachlorobenzene	0.303	0.249	17.8	86	0.00
74	Pentachlorobenzene	0.283	0.245	13.4	91	0.00
75	Carbazole	0.905	0.888	1.9	102	0.00
76	Di-n-butylphthalate	1.079	1.088	-0.8	104	0.00
77 C	Fluoranthene	1.196	1.216	-1.7	105	0.00
78 I	Chrysene-d12	1.000	1.000	0.0	116	0.00
79 S	Pyrene-d10	0.842	0.755	10.3	105	0.00
80	Pyrene	1.088	0.973	10.6	105	0.00
81	Butylbenzylphthalate	0.416	0.406	2.4	113	0.00
82	3,3'-Dichlorobenzidine	0.354	0.353	0.3	121	0.00
83	Benzo(a)anthracene	1.109	1.053	5.0	110	0.00
84	Bis(2-ethylhexyl)phthalate	0.636	0.625	1.7	115	0.00
85	Chrysene	1.037	0.967	6.8	108	0.00
86 I	Perylene-d12	1.000	1.000	0.0	120	0.00
87	Di-n-octyl phthalate	1.148	1.083	5.7	115	0.00

Data Path : Z:\HPCHEM1\BNA_M\Data\BM080315\
Data File : BM002193.D
Acq On : 03 Aug 2015 12:24
Operator : TP
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTD02094

Quant Time: Aug 04 02:31:44 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072715.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 04 02:31:26 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	Benzo(b)fluoranthene	1.124	1.050	6.6	114	0.00
89	Benzo(k)fluoranthene	1.084	1.017	6.2	113	0.00
90 S	Benzo(a)pyrene-d12	0.970	0.920	5.2	114	0.00
91 C	Benzo(a)pyrene	1.085	1.011	6.8	112	0.00
92	Indeno(1,2,3-cd)pyrene	1.250	1.112	11.0	107	0.00
93	Dibenzo(a,h)anthracene	1.070	0.947	11.5	108	0.00
94	Benzo(g,h,i)perylene	1.048	0.893	14.8	103	0.00

(#) = Out of Range

SPCC's out = 0 CCC's out = 1