

Data Path : Z:\HPCHEM1\BNA E\Data\BE100617\
 Data File : BE094194.D
 Acq On : 7 Oct 2017 00:19
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 07 02:06:57 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE100617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 06 15:16:13 2017
 Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	91	0.00
2	1,4-Dioxane	0.725	0.721	0.6	84	0.01
3	n-Nitrosodimethylamine	0.777	0.792	-1.9	93	0.01
4 S	2-Fluorophenol	1.333	1.365	-2.4	93	0.00
5 S	Phenol-d6	1.932	2.022	-4.7	96	0.00
6	bis(2-Chloroethyl)ether	1.569	1.613	-2.8	95	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	98	0.00
8 S	Nitrobenzene-d5	0.333	0.336	-0.9	100	-0.02
9	Naphthalene	4.505	4.470	0.8	96	0.00
10	Hexachlorobutadiene	0.171	0.173	-1.2	99	-0.02
11 SURR	2-Methylnaphthalene-d10	0.616	0.621	-0.8	98	0.00
12	2-Methylnaphthalene	0.740	0.750	-1.4	99	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
14 S	2,4,6-Tribromophenol	0.174	0.183	-5.2	116	0.00
15 S	2-Fluorobiphenyl	1.370	1.337	2.4	100	0.00
16	Acenaphthylene	2.026	2.022	0.2	102	0.00
17	Acenaphthene	1.338	1.317	1.6	99	0.00
18	Fluorene	1.753	1.768	-0.9	103	-0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	89	0.00
20	4-Bromophenyl-phenylether	0.230	0.266	-15.7	102	0.00
21	Hexachlorobenzene	0.207	0.259	-25.1#	112	0.00
22	Pentachlorophenol	0.056	0.065	-16.1	102	0.00
23	Phenanthrene	1.386	1.509	-8.9	96	0.00
24	Anthracene	1.057	1.102	-4.3	94	0.00
25 SURR	Fluoranthene-d10	3.995	4.364	-9.2	95	0.00
26	Fluoranthene	1.216	1.326	-9.0	96	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	94	0.00
28	Pvrene	1.376	1.376	0.0	94	0.00
29 S	Terphenyl-d14	0.755	0.761	-0.8	95	0.00
30	Benzo(a)anthracene	1.314	1.254	4.6	89	0.00
31	Chrysene	1.297	1.303	-0.5	95	0.00
32	Bis(2-ethylhexyl)phthalate	3.027	2.984	1.4	98	-0.01
33	Indeno(1,2,3-cd)pyrene	1.287	1.260	2.1	91	-0.01
34 I	Perylene-d12	1.000	1.000	0.0	93	-0.01
35	Benzo(b)fluoranthene	1.345	1.310	2.6	90	0.00
36	Benzo(k)fluoranthene	1.368	1.340	2.0	91	0.00
37 C	Benzo(a)pyrene	1.272	1.264	0.6	93	0.00
38	Dibenzo(a,h)anthracene	1.179	1.158	1.8	91	-0.01
39	Benzo(a,h,i)perylene	1.204	1.176	2.3	91	-0.01

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

(#) = Out of Range	SPCC's out = 0 CCC's out = 0			