

Data Path : Z:\HPCHEM1\BNA E\Data\BE100617\
 Data File : BE094182.D
 Acq On : 6 Oct 2017 16:51
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 07 01:32:03 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	115	0.00
2	1,4-Dioxane	0.400	0.403	-0.8	108	0.01
3	n-Nitrosodimethylamine	0.400	0.389	2.8	112	0.01
4 S	2-Fluorophenol	0.400	0.387	3.3	112	0.00
5 S	Phenol-d6	0.400	0.375	6.3	109	0.00
6	bis(2-Chloroethyl)ether	0.400	0.383	4.3	112	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	116	0.00
8 S	Nitrobenzene-d5	0.400	0.390	2.5	114	0.00
9	Naphthalene	0.400	0.402	-0.5	114	0.00
10	Hexachlorobutadiene	0.400	0.405	-1.3	117	-0.02
11 SURR	2-Methylnaphthalene-d10	0.400	0.398	0.5	114	0.00
12	2-Methylnaphthalene	0.400	0.400	0.0	115	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	119	0.00
14 S	2,4,6-Tribromophenol	0.400	0.339	15.3	108	0.00
15 S	2-Fluorobiphenyl	0.400	0.395	1.3	118	0.00
16	Acenaphthylene	0.400	0.390	2.5	116	0.00
17	Acenaphthene	0.400	0.398	0.5	116	0.00
18	Fluorene	0.400	0.388	3.0	115	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	117	0.00
20	4-Bromophenyl-phenylether	0.400	0.371	7.3	107	0.00
21	Hexachlorobenzene	0.400	0.357	10.8	105	0.00
22	Pentachlorophenol	0.400	0.319	20.3	92	0.00
23	Phenanthrene	0.400	0.377	5.8	109	0.00
24	Anthracene	0.400	0.401	-0.3	119	0.00
25 SURR	Fluoranthene-d10	0.400	0.364	9.0	105	0.00
26	Fluoranthene	0.400	0.362	9.5	105	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	105	0.00
28	Pvrene	0.400	0.394	1.5	103	0.00
29 S	Terphenyl-d14	0.400	0.390	2.5	103	0.00
30	Benzo(a)anthracene	0.400	0.388	3.0	101	0.00
31	Chrysene	0.400	0.400	0.0	105	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.345	13.8	95	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.401	-0.3	104	0.00
34 I	Perylene-d12	0.400	0.400	0.0	105	0.00
35	Benzo(b)fluoranthene	0.400	0.392	2.0	103	0.00
36	Benzo(k)fluoranthene	0.400	0.388	3.0	102	0.00
37 C	Benzo(a)pyrene	0.400	0.406	-1.5	108	0.00
38	Dibenzo(a,h)anthracene	0.400	0.389	2.8	102	0.00
39	Benzo(a,h,i)perylene	0.400	0.388	3.0	103	-0.01

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

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