

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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	91	0.00
2	1,4-Dioxane	0.400	0.398	0.5	84	0.01
3	n-Nitrosodimethylamine	0.400	0.408	-2.0	93	0.01
4 S	2-Fluorophenol	0.400	0.410	-2.5	93	0.00
5 S	Phenol-d6	0.400	0.419	-4.7	96	0.00
6	bis(2-Chloroethyl)ether	0.400	0.411	-2.7	95	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	98	0.00
8 S	Nitrobenzene-d5	0.400	0.404	-1.0	100	-0.02
9	Naphthalene	0.400	0.397	0.8	96	0.00
10	Hexachlorobutadiene	0.400	0.405	-1.3	99	-0.02
11 SURR	2-Methylnaphthalene-d10	0.400	0.403	-0.8	98	0.00
12	2-Methylnaphthalene	0.400	0.405	-1.3	99	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	103	0.00
14 S	2,4,6-Tribromophenol	0.400	0.422	-5.5	116	0.00
15 S	2-Fluorobiphenyl	0.400	0.390	2.5	100	0.00
16	Acenaphthylene	0.400	0.399	0.3	102	0.00
17	Acenaphthene	0.400	0.394	1.5	99	0.00
18	Fluorene	0.400	0.404	-1.0	103	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	89	0.00
20	4-Bromophenyl-phenylether	0.400	0.462	-15.5	102	0.00
21	Hexachlorobenzene	0.400	0.501	-25.2#	112	0.00
22	Pentachlorophenol	0.400	0.466	-16.5	102	0.00
23	Phenanthrene	0.400	0.435	-8.7	96	0.00
24	Anthracene	0.400	0.417	-4.2	94	0.00
25 SURR	Fluoranthene-d10	0.400	0.437	-9.2	95	0.00
26	Fluoranthene	0.400	0.436	-9.0	96	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	94	0.00
28	Pvrene	0.400	0.400	0.0	94	0.00
29 S	Terphenyl-d14	0.400	0.403	-0.8	95	0.00
30	Benzo(a)anthracene	0.400	0.382	4.5	89	0.00
31	Chrysene	0.400	0.402	-0.5	95	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.394	1.5	98	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.391	2.3	91	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	93	-0.01
35	Benzo(b)fluoranthene	0.400	0.389	2.8	90	0.00
36	Benzo(k)fluoranthene	0.400	0.392	2.0	91	0.00
37 C	Benzo(a)pyrene	0.400	0.398	0.5	93	0.00
38	Dibenzo(a,h)anthracene	0.400	0.393	1.8	91	-0.01
39	Benzo(a,h,i)perylene	0.400	0.390	2.5	91	-0.01

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

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