

Data Path : Z:\HPCHEM1\BNA E\DATA\BE102317\
 Data File : BE094509.D
 Acq On : 23 Oct 2017 23:26
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 24 02:22:44 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 23 22:40:47 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	98	0.00
2	1,4-Dioxane	0.400	0.445	-11.2	95	-0.01
3	n-Nitrosodimethylamine	0.400	0.411	-2.7	101	-0.01
4 S	2-Fluorophenol	0.400	0.425	-6.2	105	-0.01
5 S	Phenol-d6	0.400	0.464	-16.0	117	-0.01
6	bis(2-Chloroethyl)ether	0.400	0.402	-0.5	100	-0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	101	-0.02
8 S	Nitrobenzene-d5	0.400	0.404	-1.0	106	0.00
9	Naphthalene	0.400	0.393	1.8	97	0.00
10	Hexachlorobutadiene	0.400	0.401	-0.3	100	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.395	1.3	97	0.00
12	2-Methylnaphthalene	0.400	0.400	0.0	100	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	99	0.00
14 S	2,4,6-Tribromophenol	0.400	0.494	-23.5	127	0.00
15 S	2-Fluorobiphenyl	0.400	0.400	0.0	98	0.00
16	Acenaphthylene	0.400	0.400	0.0	100	0.00
17	Acenaphthene	0.400	0.398	0.5	97	-0.01
18	Fluorene	0.400	0.391	2.3	97	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	112	0.00
20	4-Bromophenyl-phenylether	0.400	0.412	-3.0	117	0.00
21	Hexachlorobenzene	0.400	0.376	6.0	91	0.00
22	Pentachlorophenol	0.400	0.457	-14.2	128	-0.03
23	Phenanthrene	0.400	0.340	15.0	90	0.00
24	Anthracene	0.400	0.361	9.8	98	-0.03
25 SURR	Fluoranthene-d10	0.400	0.396	1.0	102	0.00
26	Fluoranthene	0.400	0.390	2.5	101	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	99	-0.01
28	Pvrene	0.400	0.402	-0.5	101	0.00
29 S	Terphenyl-d14	0.400	0.399	0.3	101	0.00
30	Benzo(a)anthracene	0.400	0.400	0.0	99	0.00
31	Chrysene	0.400	0.404	-1.0	103	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.400	0.0	108	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.414	-3.5	104	0.00
34 I	Perylene-d12	0.400	0.400	0.0	100	0.00
35	Benzo(b)fluoranthene	0.400	0.396	1.0	101	0.00
36	Benzo(k)fluoranthene	0.400	0.393	1.8	99	0.00
37 C	Benzo(a)pyrene	0.400	0.396	1.0	99	0.00
38	Dibenzo(a,h)anthracene	0.400	0.405	-1.3	103	0.00
39	Benzo(a,h,i)perylene	0.400	0.415	-3.7	105	0.00

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

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