

Data Path : Z:\HPCHEM1\BNA E\Data\BE091817\
 Data File : BE094022.D
 Acq On : 18 Sep 2017 15:40
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Sep 19 15:30:14 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE090517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 05 13:34:56 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	129	0.00
2	1,4-Dioxane	0.400	0.445	-11.2	146	0.00
3	n-Nitrosodimethylamine	0.400	0.412	-3.0	133	0.00
4 S	2-Fluorophenol	0.400	0.388	3.0	127	-0.01
5 S	Phenol-d6	0.400	0.360	10.0	120	0.00
6	bis(2-Chloroethyl)ether	0.400	0.414	-3.5	135	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	129	0.00
8 S	Nitrobenzene-d5	0.400	0.383	4.3	124	-0.02
9	Naphthalene	0.400	0.404	-1.0	128	0.00
10	Hexachlorobutadiene	0.400	0.381	4.8	122	-0.02
11 SURR	2-Methylnaphthalene-d10	0.400	0.406	-1.5	127	0.00
12	2-Methylnaphthalene	0.400	0.402	-0.5	127	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	128	0.00
14 S	2,4,6-Tribromophenol	0.400	0.317	20.8	98	0.00
15 S	2-Fluorobiphenyl	0.400	0.397	0.8	124	0.00
16	Acenaphthylene	0.400	0.382	4.5	126	0.00
17	Acenaphthene	0.400	0.401	-0.3	126	0.00
18	Fluorene	0.400	0.389	2.8	126	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	117	-0.03
20	4-Bromophenyl-phenylether	0.400	0.384	4.0	102	0.00
21	Hexachlorobenzene	0.400	0.296	26.0#	81	0.00
22	Pentachlorophenol	0.400	0.305	23.8	89	0.00
23	Phenanthrene	0.400	0.333	16.8	95	0.00
24	Anthracene	0.400	0.403	-0.8	116	0.00
25 SURR	Fluoranthene-d10	0.400	0.373	6.8	111	0.00
26	Fluoranthene	0.400	0.372	7.0	112	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	106	-0.01
28	Pvrene	0.400	0.425	-6.2	113	0.00
29 S	Terphenyl-d14	0.400	0.414	-3.5	109	0.00
30	Benzo(a)anthracene	0.400	0.361	9.8	95	0.00
31	Chrysene	0.400	0.406	-1.5	108	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.326	18.5	90	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.272	32.0#	73	-0.01
34 I	Perylene-d12	0.400	0.400	0.0	88	0.00
35	Benzo(b)fluoranthene	0.400	0.392	2.0	86	0.00
36	Benzo(k)fluoranthene	0.400	0.463	-15.8	103	0.00
37 C	Benzo(a)pyrene	0.400	0.384	4.0	85	0.00
38	Dibenzo(a,h)anthracene	0.400	0.322	19.5	71	0.00
39	Benzo(a,h,i)perylene	0.400	0.350	12.5	78	-0.01

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

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