

Data Path : Z:\HPCHEM1\BNA E\Data\BE080317\
 Data File : BE093660.D
 Acq On : 3 Aug 2017 17:12
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Aug 04 04:56:53 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE080117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 02 10:58:31 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	109	-0.01
2	1,4-Dioxane	0.400	0.397	0.8	108	0.00
3	n-Nitrosodimethylamine	0.400	0.411	-2.7	118	0.00
4 S	2-Fluorophenol	0.400	0.442	-10.5	124	0.00
5 S	Phenol-d6	0.400	0.449	-12.2	126	-0.01
6	bis(2-Chloroethyl)ether	0.400	0.406	-1.5	113	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	116	0.00
8 S	Nitrobenzene-d5	0.400	0.410	-2.5	128	0.00
9	Naphthalene	0.400	0.389	2.8	114	0.00
10	Hexachlorobutadiene	0.400	0.387	3.3	115	-0.02
11 SURR	2-Methylnaphthalene-d10	0.400	0.400	0.0	117	0.00
12	2-Methylnaphthalene	0.400	0.397	0.8	116	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	118	0.00
14 S	2,4,6-Tribromophenol	0.400	0.439	-9.7	155	0.00
15 S	2-Fluorobiphenyl	0.400	0.389	2.8	114	-0.01
16	Acenaphthylene	0.400	0.406	-1.5	124	0.00
17	Acenaphthene	0.400	0.393	1.8	118	0.00
18	Fluorene	0.400	0.394	1.5	115	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	127	0.00
20	4-Bromophenyl-phenylether	0.400	0.373	6.8	104	0.00
21	Hexachlorobenzene	0.400	0.399	0.3	112	0.00
22	Pentachlorophenol	0.400	0.550	-37.5#	188	0.00
23	Phenanthrene	0.400	0.381	4.8	117	0.00
24	Anthracene	0.400	0.417	-4.2	125	0.00
25 SURR	Fluoranthene-d10	0.400	0.405	-1.3	118	0.00
26	Fluoranthene	0.400	0.400	0.0	117	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	116	-0.01
28	Pvrene	0.400	0.403	-0.8	117	0.00
29 S	Terphenyl-d14	0.400	0.397	0.8	113	0.00
30	Benzo(a)anthracene	0.400	0.431	-7.7	129	0.00
31	Chrysene	0.400	0.389	2.8	112	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.660	-65.0#	222	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.426	-6.5	122	0.00
34 I	Perylene-d12	0.400	0.400	0.0	123	0.00
35	Benzo(b)fluoranthene	0.400	0.380	5.0	120	0.00
36	Benzo(k)fluoranthene	0.400	0.384	4.0	122	0.00
37 C	Benzo(a)pyrene	0.400	0.401	-0.3	128	0.00
38	Dibenzo(a,h)anthracene	0.400	0.383	4.3	120	0.00
39	Benzo(a,h,i)perylene	0.400	0.382	4.5	120	0.00

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

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