

Data Path : Z:\HPCHEM1\BNA_E\Data\BE070617\
 Data File : BE093432.D
 Acq On : 6 Jul 2017 16:24
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc : LOQ 0.2 PPM
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 07 07:08:39 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE070217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 03 06:40:23 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	127	0.00
2	1,4-Dioxane	0.400	0.373	6.8	118	0.00
3	n-Nitrosodimethylamine	0.400	0.408	-2.0	158	0.00
4 S	2-Fluorophenol	0.400	0.378	5.5	123	0.00
5 S	Phenol-d6	0.400	0.410	-2.5	139	0.00
6	bis(2-Chloroethyl)ether	0.400	0.389	2.8	126	-0.01
7 I	Naphthalene-d8	0.400	0.400	0.0	142	0.00
8 S	Nitrobenzene-d5	0.400	0.434	-8.5	166	0.00
9	Naphthalene	0.400	0.386	3.5	140	0.00
10	Hexachlorobutadiene	0.400	0.377	5.8	136	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.401	-0.3	147	0.00
12	2-Methylnaphthalene	0.400	0.398	0.5	146	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	141	0.00
14 S	2,4,6-Tribromophenol	0.400	0.339	15.3	132	0.00
15 S	2-Fluorobiphenyl	0.400	0.400	0.0	146	0.00
16	Acenaphthylene	0.400	0.385	3.8	144	0.00
17	Acenaphthene	0.400	0.386	3.5	142	0.00
18	Fluorene	0.400	0.368	8.0	135	0.00
19 I	Phenanthrene-d10	0.400	0.400	0.0	121	0.00
20	4-Bromophenyl-phenylether	0.400	0.404	-1.0	132	0.00
21	Hexachlorobenzene	0.400	0.394	1.5	133	0.00
22	Pentachlorophenol	0.400	0.359	10.3	117	0.00
23	Phenanthrene	0.400	0.371	7.3	122	0.00
24	Anthracene	0.400	0.320	20.0	103	0.00
25 SURR	Fluoranthene-d10	0.400	0.330	17.5	109	0.00
26	Fluoranthene	0.400	0.333	16.8	110	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	102	0.00
28	Pyrene	0.400	0.412	-3.0	108	0.00
29 S	Terphenyl-d14	0.400	0.406	-1.5	105	0.00
30	Benzo(a)anthracene	0.400	0.386	3.5	104	0.00
31	Chrysene	0.400	0.389	2.8	101	0.00
32	Bis(2-ethylhexyl)phthalate	0.400	0.386	3.5	120	-0.01
33	Indeno(1,2,3-cd)pyrene	0.400	0.420	-5.0	110	0.00
34 I	Perylene-d12	0.400	0.400	0.0	108	0.00
35	Benzo(b)fluoranthene	0.400	0.371	7.3	103	0.00
36	Benzo(k)fluoranthene	0.400	0.391	2.3	111	0.00
37 C	Benzo(a)pyrene	0.400	0.387	3.3	111	0.00
38	Dibenzo(a,h)anthracene	0.400	0.391	2.3	110	0.00
39	Benzo(g,h,i)perylene	0.400	0.381	4.8	108	0.00

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

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