

Data Path : Z:\HPCHEM1\BNA_E\Data\BE042117\
 Data File : BE092911.D
 Acq On : 21 Apr 2017 17:03
 Operator : SJ/MA
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDC00.4

Quant Time: Apr 22 05:52:22 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE042017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 20 17:35:10 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	100	0.00
2	1,4-Dioxane	0.768	0.766	0.3	107	0.00
3	n-Nitrosodimethylamine	0.582	0.619	-6.4	112	0.00
4 S	2-Fluorophenol	0.870	0.832	4.4	103	0.00
5 S	Phenol-d6	1.213	1.190	1.9	108	0.00
6	bis(2-Chloroethyl)ether	1.474	1.498	-1.6	108	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	103	0.02
8 S	Nitrobenzene-d5	0.267	0.258	3.4	106	0.02
9	Naphthalene	1.066	1.011	5.2	103	0.00
10	Hexachlorobutadiene	0.148	0.141	4.7	102	0.00
11 SURR	2-Methylnaphthalene-d10	0.590	0.555	5.9	104	0.00
12	2-Methylnaphthalene	0.658	0.614	6.7	104	0.02
13 I	Acenaphthene-d10	1.000	1.000	0.0	106	0.00
14 S	2,4,6-Tribromophenol	0.066	0.059	10.6	110	0.01
15 S	2-Fluorobiphenyl	1.838	1.755	4.5	102	0.02
16	Acenaphthylene	6.430	5.737	10.8	102	0.00
17	Acenaphthene	1.263	1.191	5.7	103	0.00
18	Fluorene	1.541	1.422	7.7	102	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	99	0.00
20	Hexachlorobenzene	0.177	0.190	-7.3	110	0.00
21	Pentachlorophenol	0.032	0.030	6.3	112	0.00
22	Phenanthrene	1.159	1.225	-5.7	107	0.00
23	Anthracene	0.886	0.948	-7.0	113	0.00
24 SURR	Fluoranthene-d10	0.958	0.994	-3.8	111	0.00
25	Fluoranthene	4.994	4.973	0.4	107	0.00
26 I	Chrysene-d12	1.000	1.000	0.0	104	0.00
27	Pyrene	5.224	4.924	5.7	102	0.00
28 S	Terphenyl-d14	0.694	0.641	7.6	97	0.02
29	Benzo(a)anthracene	0.931	0.839	9.9	102	0.00
30	Chrysene	1.193	1.143	4.2	104	0.00
31	Bis(2-ethylhexyl)phthalate	0.241	0.227	5.8	107	0.00
32	Indeno(1,2,3-cd)pyrene	3.910	3.963	-1.4	114	0.00
33 I	Perylene-d12	1.000	1.000	0.0	130	0.01
34	Benzo(b)fluoranthene	1.401	1.063	24.1	103	0.00
35	Benzo(k)fluoranthene	1.427	1.185	17.0	117	0.01
36 C	Benzo(a)pyrene	1.133	0.885	21.9#	110	0.00
37	Dibenzo(a,h)anthracene	1.264	1.024	19.0	115	0.00
38	Benzo(g,h,i)perylene	5.484	4.491	18.1	112	0.02

(#) = Out of Range

SPCC's out = 0 CCC's out = 1