

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE051017\
 Data File : BE092979.D
 Acq On : 10 May 2017 13:21
 Operator : SJ/MA
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: May 10 14:08:30 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE051017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 10 03:43:15 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	100	0.00
2	1,4-Dioxane	0.400	0.379	5.3	103	0.00
3	n-Nitrosodimethylamine	0.400	0.398	0.5	104	-0.01
4 S	2-Fluorophenol	0.400	0.373	6.8	99	0.00
5 S	Phenol-d6	0.400	0.375	6.3	102	0.00
6	bis(2-Chloroethyl)ether	0.400	0.406	-1.5	106	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	102	0.00
8 S	Nitrobenzene-d5	0.400	0.355	11.3	97	0.00
9	Naphthalene	0.400	0.380	5.0	102	0.00
10	Hexachlorobutadiene	0.400	0.377	5.8	99	0.00
11 SURR	2-Methylnaphthalene-d10	0.400	0.384	4.0	105	0.00
12	2-Methylnaphthalene	0.400	0.386	3.5	107	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	102	0.00
14 S	2,4,6-Tribromophenol	0.400	0.322	19.5	100	0.00
15 S	2-Fluorobiphenyl	0.400	0.397	0.8	105	0.00
16	Acenaphthylene	0.400	0.362	9.5	101	0.00
17	Acenaphthene	0.400	0.383	4.3	104	0.00
18	Fluorene	0.400	0.384	4.0	102	0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	101	0.00
20	Hexachlorobenzene	0.400	0.389	2.8	106	0.00
21	Pentachlorophenol	0.400	0.371	7.3	88	0.00
22	Phenanthrene	0.400	0.417	-4.2	114	0.00
23	Anthracene	0.400	0.397	0.8	108	0.00
24 SURR	Fluoranthene-d10	0.400	0.375	6.3	105	0.00
25	Fluoranthene	0.400	0.385	3.8	108	0.00
26 I	Chrysene-d12	0.400	0.400	0.0	102	0.00
27	Pyrene	0.400	0.397	0.8	105	0.00
28 S	Terphenyl-d14	0.400	0.380	5.0	101	0.00
29	Benzo(a)anthracene	0.400	0.379	5.3	106	0.00
30	Chrysene	0.400	0.371	7.3	100	0.00
31	Bis(2-ethylhexyl)phthalate	0.400	0.411	-2.7	100	0.00
32	Indeno(1,2,3-cd)pyrene	0.400	0.362	9.5	96	0.00
33 I	Perylene-d12	0.400	0.400	0.0	100	0.00
34	Benzo(b)fluoranthene	0.400	0.350	12.5	103	0.00
35	Benzo(k)fluoranthene	0.400	0.376	6.0	99	0.00
36 C	Benzo(a)pyrene	0.400	0.344	14.0	99	0.00
37	Dibenzo(a,h)anthracene	0.400	0.352	12.0	96	0.00
38	Benzo(g,h,i)perylene	0.400	0.358	10.5	98	0.00

(#) = Out of Range

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