

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE033116\
 Data File : BE091564.D
 Acq On : 30 Mar 2016 21:22
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Mar 31 00:55:58 2016
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE033016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 30 18:45:59 2016
 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	110	0.00
2	1,4-Dioxane	0.698	0.685	1.9	104	-0.01
3	n-Nitrosodimethylamine	0.908	0.879	3.2	104	-0.03
4 S	2-Fluorophenol	1.309	1.232	5.9	103	-0.01
5 S	Phenol-d6	1.770	1.625	8.2	104	-0.02
6 I	Naphthalene-d8	1.000	1.000	0.0	109	0.00
7 S	Nitrobenzene-d5	0.336	0.311	7.4	103	-0.02
8	Naphthalene	1.033	1.014	1.8	104	-0.02
9	2-Methylnaphthalene	0.675	0.652	3.4	104	-0.02
10 I	Acenaphthene-d10	1.000	1.000	0.0	104	-0.03
11 S	2,4,6-Tribromophenol	0.222	0.163	26.6#	65	-0.01
12 S	2-Fluorobiphenyl	1.421	1.440	-1.3	103	-0.02
13	Acenaphthylene	2.016	2.043	-1.3	117	-0.01
14	Acenaphthene	1.237	1.220	1.4	103	-0.01
15	Fluorene	1.591	1.544	3.0	101	-0.01
16 I	Phenanthrene-d10	1.000	1.000	0.0	102	-0.01
17	Hexachlorobenzene	0.184	0.181	1.6	101	-0.02
18	Pentachlorophenol	0.103	0.098	4.9	111	-0.02
19	Phenanthrene	1.171	1.152	1.6	99	-0.01
20	Anthracene	1.080	1.007	6.8	97	-0.01
21	Fluoranthene	1.380	1.301	5.7	97	-0.02
22 I	Chrysene-d12	1.000	1.000	0.0	110	-0.02
23	Pyrene	1.020	0.899	11.9	96	0.00
24 S	Terphenyl-d14	0.634	0.567	10.6	96	-0.02
25	Benzo(a)anthracene	1.174	1.103	6.0	101	-0.02
26	Chrysene	1.004	0.919	8.5	98	-0.02
27	Indeno(1,2,3-cd)pyrene	1.157	1.020	11.8	98	-0.06
28 I	Perylene-d12	1.000	1.000	0.0	102	-0.03
29	Benzo(b)fluoranthene	1.213	1.208	0.4	101	-0.02
30	Benzo(k)fluoranthene	1.174	1.067	9.1	95	-0.02
31 C	Benzo(a)pyrene	1.108	1.036	6.5	97	-0.03
32	Dibenzo(a,h)anthracene	1.088	1.027	5.6	97	-0.05
33	Benzo(g,h,i)perylene	1.111	1.051	5.4	97	-0.05

(#) = Out of Range

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