

Data Path : Z:\HPCHEM1\BNA E\DATA\BE100716\
 Data File : BE092443.D
 Acq On : 7 Oct 2016 11:53
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 07 23:06:54 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE092216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 22 18:25:56 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	109	-0.02
2	1,4-Dioxane	0.635	0.661	-4.1	113	-0.01
3	n-Nitrosodimethylamine	0.588	0.503	14.5	103	0.00
4 S	2-Fluorophenol	0.999	0.825	17.4	97	0.00
5 S	Phenol-d6	1.388	1.000	28.0#	91	0.00
6	bis(2-Chloroethyl)ether	1.222	0.857	29.9#	98	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	104	0.00
8 S	Nitrobenzene-d5	0.299	0.244	18.4	101	0.00
9	Naphthalene	1.082	1.076	0.6	104	-0.02
10	Hexachlorobutadiene	0.170	0.169	0.6	104	-0.02
11	2-Methylnaphthalene	0.716	0.698	2.5	103	0.00
12 I	Acenaphthene-d10	1.000	1.000	0.0	102	0.00
13 S	2,4,6-Tribromophenol	0.144	0.101	29.9#	83	0.00
14 S	2-Fluorobiphenyl	1.184	1.213	-2.4	106	0.00
15	Acenaphthylene	7.240	6.898	4.7	100	0.00
16	Acenaphthene	1.134	1.135	-0.1	99	0.00
17	Fluorene	1.438	1.382	3.9	99	0.00
18 I	Phenanthrene-d10	1.000	1.000	0.0	93	0.02
19	Hexachlorobenzene	0.214	0.229	-7.0	106	0.00
20	Pentachlorophenol	0.050	0.054	-8.0	119	0.00
21	Phenanthrene	1.149	1.242	-8.1	106	0.00
22	Anthracene	1.105	1.053	4.7	97	0.02
23	Fluoranthene	5.460	5.122	6.2	94	0.00
24 I	Chrysene-d12	1.000	1.000	0.0	94	0.00
25	Pyrene	4.333	4.259	1.7	96	0.00
26 S	Terphenyl-d14	0.620	0.625	-0.8	95	0.02
27	Benzo(a)anthracene	4.873	4.475	8.2	88	0.00
28	Chrysene	4.462	4.853	-8.8	95	0.00
29	Bis(2-ethylhexyl)phthalate	0.420	0.325	22.6	72	0.00
30	Indeno(1,2,3-cd)pyrene	4.734	4.018	15.1	81	0.03
31 I	Perylene-d12	1.000	1.000	0.0	88	0.02
32	Benzo(b)fluoranthene	1.211	1.129	6.8	82	0.02
33	Benzo(k)fluoranthene	1.267	1.230	2.9	92	0.02
34 C	Benzo(a)pyrene	1.187	1.077	9.3	85	0.02
35	Dibenzo(a,h)anthracene	1.088	0.956	12.1	81	0.05
36	Benzo(g,h,i)perylene	4.673	4.337	7.2	84	0.03

(#) = Out of Range

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