

Data Path : Z:\HPCHEM1\BNA E\DATA\BE111716\
 Data File : BE092561.D
 Acq On : 17 Nov 2016 16:14
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Nov 18 00:19:14 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE111616.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 16 15:43:18 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	111	-0.02
2	1,4-Dioxane	0.681	0.821	-20.6	140	0.00
3	n-Nitrosodimethylamine	0.654	0.633	3.2	126	-0.01
4 S	2-Fluorophenol	0.844	0.907	-7.5	120	0.00
5 S	Phenol-d6	1.104	1.158	-4.9	127	0.00
6	bis(2-Chloroethyl)ether	1.341	1.193	11.0	112	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	110	0.00
8 S	Nitrobenzene-d5	0.269	0.271	-0.7	113	-0.02
9	Naphthalene	1.032	1.038	-0.6	111	-0.02
10	Hexachlorobutadiene	0.157	0.160	-1.9	111	0.00
11	2-Methylnaphthalene	0.658	0.654	0.6	115	-0.01
12 I	Acenaphthene-d10	1.000	1.000	0.0	115	0.00
13 S	2,4,6-Tribromophenol	0.112	0.112	0.0	122	0.00
14 S	2-Fluorobiphenyl	1.219	1.210	0.7	113	0.00
15	Acenaphthylene	6.962	6.847	1.7	118	0.00
16	Acenaphthene	1.136	1.119	1.5	113	0.00
17	Fluorene	1.408	1.392	1.1	115	-0.01
18 I	Phenanthrene-d10	1.000	1.000	0.0	105	0.00
19	Hexachlorobenzene	0.197	0.237	-20.3	140	0.00
20	Pentachlorophenol	0.036	0.040	-11.1	149	0.00
21	Phenanthrene	1.119	1.175	-5.0	114	-0.02
22	Anthracene	1.053	1.060	-0.7	112	0.00
23	Fluoranthene	5.219	5.307	-1.7	113	0.00
24 I	Chrysene-d12	1.000	1.000	0.0	108	0.00
25	Pyrene	3.975	3.925	1.3	114	0.00
26 S	Terphenyl-d14	0.572	0.556	2.8	113	-0.02
27	Benzo(a)anthracene	4.692	4.342	7.5	119	0.00
28	Chrysene	4.250	4.509	-6.1	110	-0.02
29	Bis(2-ethylhexyl)phthalate	0.305	0.355	-16.4	150#	-0.02
30	Indeno(1,2,3-cd)pyrene	4.722	4.284	9.3	103	0.00
31 I	Perylene-d12	1.000	1.000	0.0	102	-0.01
32	Benzo(b)fluoranthene	1.213	1.152	5.0	101	-0.01
33	Benzo(k)fluoranthene	1.270	1.263	0.6	107	-0.01
34 C	Benzo(a)pyrene	1.122	1.077	4.0	106	-0.01
35	Dibenzo(a,h)anthracene	1.111	1.044	6.0	103	-0.02
36	Benzo(g,h,i)perylene	4.702	4.497	4.4	102	0.00

(#) = Out of Range

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