

Data Path : Z:\HPCHEM1\BNA E\DATA\BE092316\
 Data File : BE092423.D
 Acq On : 23 Sep 2016 19:50
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Sep 23 23:17:43 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	98	-0.02
2	1,4-Dioxane	0.635	0.628	1.1	96	-0.01
3	n-Nitrosodimethylamine	0.588	0.445	24.3	83	0.00
4 S	2-Fluorophenol	0.999	0.748	25.1#	79	0.00
5 S	Phenol-d6	1.388	0.991	28.6#	81	0.00
6	bis(2-Chloroethyl)ether	1.222	0.790	35.4#	82	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	95	0.00
8 S	Nitrobenzene-d5	0.299	0.259	13.4	97	0.00
9	Naphthalene	1.082	1.042	3.7	92	0.00
10	Hexachlorobutadiene	0.170	0.164	3.5	92	0.00
11	2-Methylnaphthalene	0.716	0.679	5.2	91	0.00
12 I	Acenaphthene-d10	1.000	1.000	0.0	96	0.00
13 S	2,4,6-Tribromophenol	0.144	0.100	30.6#	77	0.01
14 S	2-Fluorobiphenyl	1.184	1.136	4.1	93	0.01
15	Acenaphthylene	7.240	6.651	8.1	90	0.02
16	Acenaphthene	1.134	1.105	2.6	90	0.00
17	Fluorene	1.438	1.327	7.7	89	0.00
18 I	Phenanthrene-d10	1.000	1.000	0.0	90	0.02
19	Hexachlorobenzene	0.214	0.206	3.7	92	0.00
20	Pentachlorophenol	0.050	0.029	42.0#	62	0.00
21	Phenanthrene	1.149	1.099	4.4	90	0.02
22	Anthracene	1.105	0.984	11.0	88	0.02
23	Fluoranthene	5.460	4.898	10.3	87	0.02
24 I	Chrysene-d12	1.000	1.000	0.0	82	0.00
25	Pyrene	4.333	4.327	0.1	86	0.02
26 S	Terphenyl-d14	0.620	0.636	-2.6	85	0.02
27	Benzo(a)anthracene	4.873	4.228	13.2	73	0.00
28	Chrysene	4.462	4.938	-10.7	85	0.02
29	Bis(2-ethylhexyl)phthalate	0.420	0.324	22.9	63	0.02
30	Indeno(1,2,3-cd)pyrene	4.734	4.245	10.3	75	0.07
31 I	Perylene-d12	1.000	1.000	0.0	88	0.05
32	Benzo(b)fluoranthene	1.211	1.049	13.4	76	0.03
33	Benzo(k)fluoranthene	1.267	1.188	6.2	88	0.03
34 C	Benzo(a)pyrene	1.187	1.054	11.2	82	0.03
35	Dibenzo(a,h)anthracene	1.088	0.928	14.7	78	0.07
36	Benzo(g,h,i)perylene	4.673	4.147	11.3	80	0.07

(#) = Out of Range

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