

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042016\
 Data File : BE091694.D
 Acq On : 20 Apr 2016 20:47
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Apr 21 04:24:29 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE042016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 20 18:55:38 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	0.00
2	1,4-Dioxane	0.615	0.669	-8.8	95	0.00
3	n-Nitrosodimethylamine	0.718	0.706	1.7	101	0.00
4 S	2-Fluorophenol	1.106	1.055	4.6	97	0.00
5 S	Phenol-d6	1.496	1.380	7.8	96	0.00
6	bis(2-Chloroethyl)ether	1.344	1.272	5.4	95	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	100	0.00
8 S	Nitrobenzene-d5	0.267	0.243	9.0	102	0.00
9	Nitrobenzene	0.267	0.247	7.5	102	-0.01
10	Naphthalene	1.019	1.002	1.7	97	0.00
11	Hexachlorobutadiene	0.146	0.147	-0.7	100	0.00
12	2-Methylnaphthalene	0.654	0.634	3.1	98	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	98	0.00
14 S	2,4,6-Tribromophenol	0.117	0.100	14.5	106	-0.01
15 S	2-Fluorobiphenyl	1.383	1.393	-0.7	99	0.00
16	Acenaphthylene	1.999	1.849	7.5	100	0.00
17	Acenaphthene	1.231	1.209	1.8	98	0.00
18	Fluorene	1.526	1.508	1.2	99	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	100	0.00
20	4-Bromophenyl-phenylether	0.171	0.169	1.2	100	0.00
21	Hexachlorobenzene	0.183	0.183	0.0	98	0.00
22	Pentachlorophenol	0.060	0.040	33.3#	96	-0.02
23	Phenanthrene	1.123	1.161	-3.4	99	-0.01
24	Anthracene	1.002	0.956	4.6	100	-0.01
25	Fluoranthene	1.110	1.020	8.1	95	-0.02
26 I	Chrysene-d12	1.000	1.000	0.0	94	0.00
27	Benzydine	0.292	0.188	35.6#	99	0.00
28	Pvrene	1.155	1.141	1.2	99	0.00
29 S	Terphenyl-d14	0.793	0.769	3.0	97	0.00
30	Benzo(a)anthracene	1.215	1.124	7.5	94	0.00
31	3,3'-Dichlorobenzidine	0.510	0.404	20.8	103	-0.02
32	Chrysene	1.324	1.429	-7.9	104	0.00
33	Bis(2-ethylhexyl)phthalate	0.233	0.204	12.4	113	-0.02
34	Indeno(1,2,3-cd)pyrene	1.185	1.084	8.5	94	-0.01
35 I	Perylene-d12	1.000	1.000	0.0	95	-0.01
36	Benzo(b)fluoranthene	1.199	1.126	6.1	94	0.00
37	Benzo(k)fluoranthene	1.180	1.106	6.3	94	-0.01
38 C	Benzo(a)pyrene	1.037	0.933	10.0	93	-0.01
39	Dibenzo(a,h)anthracene	1.063	0.954	10.3	93	0.00
40	Benzo(a,h,i)perylene	1.090	1.013	7.1	94	-0.01

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

(#) = Out of Range				
SPCC's out = 0 CCC's out = 0				