

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042616\
 Data File : BE091710.D
 Acq On : 26 Apr 2016 18:43
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Apr 27 01:04:29 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE042616.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 26 16:09:46 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	95	0.00
2	1,4-Dioxane	0.752	0.656	12.8	82	0.02
3	n-Nitrosodimethylamine	0.908	0.827	8.9	88	0.02
4 S	2-Fluorophenol	1.228	1.151	6.3	92	0.00
5 S	Phenol-d6	1.629	1.466	10.0	91	0.00
6	bis(2-Chloroethyl)ether	1.433	1.414	1.3	93	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	93	0.00
8 S	Nitrobenzene-d5	0.323	0.296	8.4	91	0.00
9	Nitrobenzene	0.338	0.303	10.4	90	0.00
10	Naphthalene	1.010	1.034	-2.4	94	0.00
11	Hexachlorobutadiene	0.149	0.152	-2.0	97	0.00
12	2-Methylnaphthalene	0.649	0.627	3.4	90	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	92	0.00
14 S	2,4,6-Tribromophenol	0.157	0.134	14.6	86	0.00
15 S	2-Fluorobiphenyl	1.414	1.394	1.4	91	0.00
16	Acenaphthylene	1.923	1.892	1.6	102	0.00
17	Acenaphthene	1.247	1.220	2.2	90	0.00
18	Fluorene	1.550	1.522	1.8	90	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	90	0.00
20	4-Bromophenyl-phenylether	0.180	0.172	4.4	85	0.00
21	Hexachlorobenzene	0.180	0.177	1.7	89	0.00
22	Pentachlorophenol	0.085	0.065	23.5	81	0.00
23	Phenanthrene	1.136	1.166	-2.6	91	0.00
24	Anthracene	1.025	0.981	4.3	88	0.00
25	Fluoranthene	1.188	1.181	0.6	89	0.00
26 I	Chrysene-d12	1.000	1.000	0.0	81	0.00
27	Benzidine	0.286	0.266	7.0	93	0.00
28	Pvrene	1.130	1.213	-7.3	90	0.00
29 S	Terphenyl-d14	0.780	0.837	-7.3	89	0.00
30	Benzo(a)anthracene	1.221	1.267	-3.8	86	0.00
31	3,3'-Dichlorobenzidine	0.623	0.614	1.4	92	-0.02
32	Chrysene	1.269	1.554	-22.5	94	0.00
33	Bis(2-ethylhexyl)phthalate	0.370	0.365	1.4	104	-0.02
34	Indeno(1,2,3-cd)pyrene	1.130	1.273	-12.7	90	0.00
35 I	Perylene-d12	1.000	1.000	0.0	91	0.00
36	Benzo(b)fluoranthene	1.172	1.183	-0.9	94	0.00
37	Benzo(k)fluoranthene	1.191	1.080	9.3	86	0.00
38 C	Benzo(a)pyrene	1.103	1.034	6.3	90	0.00
39	Dibenzo(a,h)anthracene	1.045	0.997	4.6	89	-0.01
40	Benzo(a,h,i)perylene	1.062	1.009	5.0	89	-0.01

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(#) = Out of Range

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