

Data Path : Z:\HPCHEM1\BNA E\DATA\BE041816\
 Data File : BE091669.D
 Acq On : 18 Apr 2016 17:45
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Apr 19 05:20:45 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE041816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 18 17:20:40 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	90	0.00
2	1,4-Dioxane	0.645	0.648	-0.5	91	0.00
3	n-Nitrosodimethylamine	0.717	0.737	-2.8	100	0.00
4 S	2-Fluorophenol	1.088	1.127	-3.6	99	0.00
5 S	Phenol-d6	1.467	1.450	1.2	97	0.00
6	bis(2-Chloroethyl)ether	1.348	1.361	-1.0	95	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	93	0.00
8 S	Nitrobenzene-d5	0.256	0.251	2.0	106	0.00
9	Nitrobenzene	0.258	0.254	1.6	104	0.00
10	Naphthalene	1.018	1.036	-1.8	95	0.00
11	Hexachlorobutadiene	0.148	0.148	0.0	95	0.00
12	2-Methylnaphthalene	0.650	0.656	-0.9	97	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	95	0.00
14 S	2,4,6-Tribromophenol	0.106	0.103	2.8	113	-0.01
15 S	2-Fluorobiphenyl	1.393	1.408	-1.1	97	0.00
16	Acenaphthylene	1.840	1.859	-1.0	110	0.00
17	Acenaphthene	1.230	1.285	-4.5	96	0.00
18	Fluorene	1.507	1.508	-0.1	97	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	86	0.00
20	4-Bromophenyl-phenylether	0.173	0.168	2.9	97	0.00
21	Hexachlorobenzene	0.184	0.182	1.1	94	0.00
22	Pentachlorophenol	0.061	0.054	11.5	128	0.00
23	Phenanthrene	1.136	1.171	-3.1	97	0.00
24	Anthracene	1.009	0.989	2.0	98	0.00
25	Fluoranthene	1.129	1.147	-1.6	101	0.00
26 I	Chrysene-d12	1.000	1.000	0.0	95	0.00
27	Benzydine	0.363	0.278	23.4	140	-0.02
28	Pvrene	0.999	1.171	-17.2	118	0.00
29 S	Terphenyl-d14	0.681	0.758	-11.3	110	0.00
30	Benzo(a)anthracene	1.160	1.236	-6.6	106	0.00
31	3,3'-Dichlorobenzidine	0.454	0.471	-3.7	128	0.00
32	Chrysene	1.176	1.253	-6.5	100	0.00
33	Bis(2-ethylhexyl)phthalate	0.212	0.273	-28.8#	149	0.00
34	Indeno(1,2,3-cd)pyrene	1.070	1.128	-5.4	103	0.00
35 I	Perylene-d12	1.000	1.000	0.0	97	-0.01
36	Benzo(b)fluoranthene	1.100	1.158	-5.3	109	0.00
37	Benzo(k)fluoranthene	1.168	1.209	-3.5	105	0.00
38 C	Benzo(a)pyrene	1.026	1.078	-5.1	110	0.00
39	Dibenzo(a,h)anthracene	1.003	1.012	-0.9	104	0.00
40	Benzo(a,h,i)perylene	1.039	1.041	-0.2	102	0.00

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)
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(#) = Out of Range

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