

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042916\
 Data File : BE091739.D
 Acq On : 29 Apr 2016 19:07
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Apr 29 23:55:12 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	93	0.00
2	1,4-Dioxane	0.752	0.642	14.6	78	0.02
3	n-Nitrosodimethylamine	0.908	0.746	17.8	78	0.02
4 S	2-Fluorophenol	1.228	1.069	12.9	83	0.00
5 S	Phenol-d6	1.629	1.336	18.0	81	0.02
6	bis(2-Chloroethyl)ether	1.433	1.347	6.0	86	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	91	0.00
8 S	Nitrobenzene-d5	0.323	0.264	18.3	80	0.01
9	Nitrobenzene	0.338	0.268	20.7	79	0.00
10	Naphthalene	1.010	1.028	-1.8	91	0.00
11	Hexachlorobutadiene	0.149	0.151	-1.3	94	-0.02
12	2-Methylnaphthalene	0.649	0.620	4.5	87	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	87	-0.01
14 S	2,4,6-Tribromophenol	0.157	0.117	25.5#	71	0.00
15 S	2-Fluorobiphenyl	1.414	1.411	0.2	88	0.00
16	Acenaphthylene	1.923	1.838	4.4	94	0.00
17	Acenaphthene	1.247	1.199	3.8	84	0.00
18	Fluorene	1.550	1.515	2.3	85	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	88	0.00
20	4-Bromophenyl-phenylether	0.180	0.167	7.2	80	0.00
21	Hexachlorobenzene	0.180	0.180	0.0	88	0.00
22	Pentachlorophenol	0.085	0.053	37.6#	64	0.00
23	Phenanthrene	1.136	1.152	-1.4	87	0.00
24	Anthracene	1.025	0.962	6.1	84	0.00
25	Fluoranthene	1.188	1.202	-1.2	88	0.00
26 I	Chrysene-d12	1.000	1.000	0.0	78	0.00
27	Benzydine	0.286	0.215	24.8	72	0.00
28	Pvrene	1.130	1.220	-8.0	87	0.00
29 S	Terphenyl-d14	0.780	0.821	-5.3	84	0.00
30	Benzo(a)anthracene	1.221	1.198	1.9	78	0.00
31	3,3'-Dichlorobenzidine	0.623	0.525	15.7	75	-0.02
32	Chrysene	1.269	1.476	-16.3	85	0.00
33	Bis(2-ethylhexyl)phthalate	0.370	0.356	3.8	97	-0.02
34	Indeno(1,2,3-cd)pyrene	1.130	1.291	-14.2	87	-0.01
35 I	Perylene-d12	1.000	1.000	0.0	88	-0.01
36	Benzo(b)fluoranthene	1.172	1.234	-5.3	96	-0.01
37	Benzo(k)fluoranthene	1.191	1.037	12.9	80	0.00
38 C	Benzo(a)pyrene	1.103	1.036	6.1	87	0.00
39	Dibenzo(a,h)anthracene	1.045	0.995	4.8	86	-0.02
40	Benzo(a,h,i)perylene	1.062	0.999	5.9	86	-0.02

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Compound	AvgRF	CCRF	%Dev Area	% Dev(min)

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