

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE101717\
 Data File : BE094435.D
 Acq On : 17 Oct 2017 18:39
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 17 20:35:49 2017
 Quant Method : Z:\HPCHEM1\BNA_E\Methods\8270-SIM-BE101617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 14:54:40 2017
 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	96	-0.02
2	1,4-Dioxane	0.698	0.721	-3.3	98	0.00
3	n-Nitrosodimethylamine	0.695	0.600	13.7	83	0.00
4 S	2-Fluorophenol	1.352	1.307	3.3	91	0.00
5 S	Phenol-d6	2.108	2.021	4.1	89	-0.02
6	bis(2-Chloroethyl)ether	1.559	1.532	1.7	91	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	95	0.00
8 S	Nitrobenzene-d5	0.348	0.333	4.3	92	0.00
9	Naphthalene	4.473	4.460	0.3	95	-0.02
10	Hexachlorobutadiene	0.179	0.170	5.0	91	-0.02
11 SURR	2-Methylnaphthalene-d10	0.617	0.627	-1.6	97	0.00
12	2-Methylnaphthalene	0.737	0.757	-2.7	99	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	105	0.00
14 S	2,4,6-Tribromophenol	0.152	0.138	9.2	90	0.00
15 S	2-Fluorobiphenyl	1.428	1.375	3.7	99	0.00
16	Acenaphthylene	2.169	2.130	1.8	103	0.00
17	Acenaphthene	1.366	1.331	2.6	103	0.00
18	Fluorene	1.761	1.749	0.7	105	-0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	98	-0.03
20	4-Bromophenyl-phenylether	0.200	0.209	-4.5	114	-0.03
21	Hexachlorobenzene	0.365	0.347	4.9	96	0.00
22	Pentachlorophenol	0.032	0.023	28.1#	78	0.03
23	Phenanthrene	1.574	1.373	12.8	82	0.00
24	Anthracene	1.216	1.241	-2.1	103	0.00
25 SURR	Fluoranthene-d10	6.441	6.411	0.5	101	0.00
26	Fluoranthene	1.954	1.954	0.0	103	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	109	0.00
28	Pyrene	1.509	1.391	7.8	102	0.00
29 S	Terphenyl-d14	0.807	0.754	6.6	101	0.00
30	Benzo(a)anthracene	1.384	1.334	3.6	106	-0.01
31	Chrysene	1.332	1.270	4.7	102	-0.01
32	Bis(2-ethylhexyl)phthalate	3.984	3.572	10.3	102	0.00
33	Indeno(1,2,3-cd)pyrene	1.341	1.287	4.0	105	-0.02
34 I	Perylene-d12	1.000	1.000	0.0	109	-0.02
35	Benzo(b)fluoranthene	1.357	1.286	5.2	103	0.00
36	Benzo(k)fluoranthene	1.337	1.350	-1.0	109	0.00
37 C	Benzo(a)pyrene	1.275	1.233	3.3	105	-0.01
38	Dibenzo(a,h)anthracene	1.207	1.158	4.1	103	-0.01
39	Benzo(g,h,i)perylene	1.161	1.130	2.7	105	-0.02

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