

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE051718\
 Data File : BE096615.D
 Acq On : 17 May 2018 18:30
 Operator : SJ/JU
 Sample : SSTDC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDC0.4

Quant Time: May 18 05:53:25 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE051518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 15 13:02:03 2018
 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	106	0.00
2	1,4-Dioxane	0.600	0.567	5.5	112	0.00
3	n-Nitrosodimethylamine	0.750	0.698	6.9	108	0.00
4 S	2-Fluorophenol	1.103	1.000	9.3	106	0.00
5 S	Phenol-d6	1.561	1.404	10.1	101	0.00
6	bis(2-Chloroethyl)ether	1.460	1.405	3.8	108	-0.01
7 I	Naphthalene-d8	1.000	1.000	0.0	107	-0.01
8 S	Nitrobenzene-d5	0.426	0.408	4.2	110	0.00
9	Naphthalene	4.208	3.930	6.6	106	0.00
10	Hexachlorobutadiene	0.235	0.212	9.8	111	-0.01
11 SURR	2-Methylnaphthalene-d10	0.653	0.597	8.6	104	0.00
12	2-Methylnaphthalene	0.721	0.662	8.2	105	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	99	0.00
14 S	2,4,6-Tribromophenol	0.158	0.144	8.9	98	0.00
15 S	2-Fluorobiphenyl	1.717	1.703	0.8	106	0.00
16	Acenaphthylene	2.088	2.005	4.0	104	0.00
17	Acenaphthene	1.272	1.216	4.4	101	-0.01
18	Fluorene	1.583	1.487	6.1	100	-0.01
19 I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
20	4-Bromophenyl-phenylether	0.243	0.235	3.3	101	0.00
21	Hexachlorobenzene	0.253	0.244	3.6	101	0.00
22	Pentachlorophenol	0.053	0.038	28.3#	109	0.00
23	Phenanthrene	1.123	1.063	5.3	98	0.00
24	Anthracene	1.045	0.963	7.8	96	0.00
25 SURR	Fluoranthene-d10	4.890	4.508	7.8	97	0.00
26	Fluoranthene	1.343	1.251	6.9	98	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	98	0.00
28	Pyrene	0.722	0.707	2.1	102	0.00
29 S	Terphenyl-d14	0.905	0.855	5.5	96	0.00
30	Benzo(a)anthracene	1.230	1.136	7.6	97	0.00
31	Chrysene	1.280	1.217	4.9	99	0.00
32	Bis(2-ethylhexyl)phthalate	2.632	2.285	13.2	95	0.00
33	Indeno(1,2,3-cd)pyrene	1.227	1.154	5.9	97	0.00
34 I	Perylene-d12	1.000	1.000	0.0	98	0.00
35	Benzo(b)fluoranthene	1.234	1.173	4.9	100	0.00
36	Benzo(k)fluoranthene	1.306	1.244	4.7	100	0.00
37 C	Benzo(a)pyrene	1.180	1.114	5.6	98	0.00
38	Dibenzo(a,h)anthracene	1.079	0.988	8.4	95	0.00
39	Benzo(g,h,i)perylene	1.129	1.067	5.5	99	0.00

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