

Data Path : Z:\HPCHEM1\BNA_E\Data\BE070117\
 Data File : BE093403.D
 Acq On : 1 Jul 2017 22:22
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 03 06:05:42 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE070117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 01 15:43:24 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	103	0.00
2	1,4-Dioxane	0.826	0.774	6.3	99	0.00
3	n-Nitrosodimethylamine	0.497	0.429	13.7	106	0.00
4 S	2-Fluorophenol	1.247	1.189	4.7	101	0.00
5 S	Phenol-d6	1.861	1.756	5.6	104	0.00
6	bis(2-Chloroethyl)ether	1.312	1.279	2.5	109	0.00
7 I	Naphthalene-d8	1.000	1.000	0.0	103	0.00
8 S	Nitrobenzene-d5	0.238	0.260	-9.2	129	-0.02
9	Naphthalene	4.155	4.038	2.8	104	-0.02
10	Hexachlorobutadiene	0.167	0.166	0.6	103	0.00
11 SURR	2-Methylnaphthalene-d10	0.639	0.621	2.8	104	0.00
12	2-Methylnaphthalene	0.708	0.687	3.0	103	0.00
13 I	Acenaphthene-d10	1.000	1.000	0.0	106	-0.02
14 S	2,4,6-Tribromophenol	0.172	0.168	2.3	114	0.00
15 S	2-Fluorobiphenyl	1.566	1.499	4.3	103	0.00
16	Acenaphthylene	1.739	1.556	10.5	100	-0.02
17	Acenaphthene	1.224	1.194	2.5	107	-0.02
18	Fluorene	1.683	1.648	2.1	105	0.00
19 I	Phenanthrene-d10	1.000	1.000	0.0	88	0.00
20	4-Bromophenyl-phenylether	0.091	0.170	-86.8#	162#	0.00
21	Hexachlorobenzene	0.153	0.221	-44.4#	125	0.00
22	Pentachlorophenol	0.018	0.027	-50.0#	158#	-0.03
23	Phenanthrene	0.968	1.251	-29.2#	109	0.00
24	Anthracene	0.974	0.828	15.0	81	0.00
25 SURR	Fluoranthene-d10	3.688	3.941	-6.9	99	0.00
26	Fluoranthene	1.049	1.120	-6.8	99	0.00
27 I	Chrysene-d12	1.000	1.000	0.0	97	-0.01
28	Pyrene	1.056	1.060	-0.4	100	0.00
29 S	Terphenyl-d14	0.742	0.742	0.0	99	0.00
30	Benzo(a)anthracene	1.088	1.045	4.0	98	-0.01
31	Chrysene	1.137	1.135	0.2	99	-0.01
32	Bis(2-ethylhexyl)phthalate	1.612	1.408	12.7	102	-0.01
33	Indeno(1,2,3-cd)pyrene	1.112	1.088	2.2	98	-0.02
34 I	Perylene-d12	1.000	1.000	0.0	98	0.00
35	Benzo(b)fluoranthene	1.256	1.214	3.3	98	0.00
36	Benzo(k)fluoranthene	1.260	1.217	3.4	99	-0.01
37 C	Benzo(a)pyrene	1.116	1.057	5.3	98	-0.01
38	Dibenzo(a,h)anthracene	1.137	1.099	3.3	99	-0.02
39	Benzo(g,h,i)perylene	1.159	1.120	3.4	97	-0.02

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