

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE121517\
 Data File : BE095002.D
 Acq On : 15 Dec 2017 15:15
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Dec 15 16:32:54 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 14 18:42:53 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	94	0.00
2	1,4-Dioxane	0.400	0.383	4.3	85	0.00
3	n-Nitrosodimethylamine	0.400	0.404	-1.0	95	0.00
4 S	2-Fluorophenol	0.400	0.420	-5.0	103	0.00
5 S	Phenol-d6	0.400	0.425	-6.2	103	0.00
6	bis(2-Chloroethyl)ether	0.400	0.400	0.0	95	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	100	-0.02
8 S	Nitrobenzene-d5	0.400	0.446	-11.5	117	0.00
9	Naphthalene	0.400	0.390	2.5	98	-0.02
10	Hexachlorobutadiene	0.400	0.382	4.5	95	-0.01
11 SURR	2-Methylnaphthalene-d10	0.400	0.390	2.5	99	-0.01
12	2-Methylnaphthalene	0.400	0.362	9.5	92	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	100	-0.02
14 S	2,4,6-Tribromophenol	0.400	0.474	-18.5	127	-0.02
15 S	2-Fluorobiphenyl	0.400	0.390	2.5	96	0.00
16	Acenaphthylene	0.400	0.388	3.0	102	-0.02
17	Acenaphthene	0.400	0.392	2.0	99	-0.02
18	Fluorene	0.400	0.384	4.0	98	-0.02
19 I	Phenanthrene-d10	0.400	0.400	0.0	100	-0.02
20	4-Bromophenyl-phenylether	0.400	0.387	3.3	100	-0.02
21	Hexachlorobenzene	0.400	0.378	5.5	96	-0.02
22	Pentachlorophenol	0.400	0.480	-20.0	135	-0.02
23	Phenanthrene	0.400	0.387	3.3	100	-0.02
24	Anthracene	0.400	0.381	4.8	101	-0.02
25 SURR	Fluoranthene-d10	0.400	0.381	4.8	100	-0.01
26	Fluoranthene	0.400	0.379	5.3	98	-0.01
27 I	Chrysene-d12	0.400	0.400	0.0	103	-0.02
28	Pyrene	0.400	0.383	4.3	96	-0.02
29 S	Terphenyl-d14	0.400	0.382	4.5	98	-0.02
30	Benzo(a)anthracene	0.400	0.382	4.5	100	-0.02
31	Chrysene	0.400	0.375	6.3	97	-0.02
32	Bis(2-ethylhexyl)phthalate	0.400	0.433	-8.2	125	-0.03
33	Indeno(1,2,3-cd)pyrene	0.400	0.393	1.8	99	-0.05
34 I	Perylene-d12	0.400	0.400	0.0	99	-0.03
35	Benzo(b)fluoranthene	0.400	0.380	5.0	94	-0.02
36	Benzo(k)fluoranthene	0.400	0.377	5.8	98	-0.03
37 C	Benzo(a)pyrene	0.400	0.388	3.0	97	-0.02
38	Dibenzo(a,h)anthracene	0.400	0.385	3.8	97	-0.04
39	Benzo(g,h,i)perylene	0.400	0.387	3.3	98	-0.05

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

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