

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE101917\
 Data File : BE094457.D
 Acq On : 19 Oct 2017 21:30
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Oct 20 13:52:26 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	0.400	0.400	0.0	94	-0.01
2	1,4-Dioxane	0.400	0.422	-5.5	96	0.00
3	n-Nitrosodimethylamine	0.400	0.385	3.8	90	0.01
4 S	2-Fluorophenol	0.400	0.427	-6.7	99	0.00
5 S	Phenol-d6	0.400	0.420	-5.0	95	0.00
6	bis(2-Chloroethyl)ether	0.400	0.427	-6.7	96	0.00
7 I	Naphthalene-d8	0.400	0.400	0.0	91	-0.02
8 S	Nitrobenzene-d5	0.400	0.413	-3.2	95	0.00
9	Naphthalene	0.400	0.438	-9.5	100	-0.02
10	Hexachlorobutadiene	0.400	0.417	-4.2	96	-0.02
11 SURR	2-Methylnaphthalene-d10	0.400	0.446	-11.5	101	0.00
12	2-Methylnaphthalene	0.400	0.450	-12.5	104	0.00
13 I	Acenaphthene-d10	0.400	0.400	0.0	100	-0.01
14 S	2,4,6-Tribromophenol	0.400	0.393	1.8	92	0.00
15 S	2-Fluorobiphenyl	0.400	0.427	-6.7	104	0.00
16	Acenaphthylene	0.400	0.420	-5.0	104	0.00
17	Acenaphthene	0.400	0.420	-5.0	105	0.00
18	Fluorene	0.400	0.416	-4.0	104	-0.01
19 I	Phenanthrene-d10	0.400	0.400	0.0	101	-0.03
20	4-Bromophenyl-phenylether	0.400	0.460	-15.0	130	-0.03
21	Hexachlorobenzene	0.400	0.296	26.0#	81	0.00
22	Pentachlorophenol	0.400	0.470	-17.5	113	0.00
23	Phenanthrene	0.400	0.308	23.0	75	0.00
24	Anthracene	0.400	0.407	-1.7	107	-0.03
25 SURR	Fluoranthene-d10	0.400	0.373	6.8	98	0.00
26	Fluoranthene	0.400	0.372	7.0	99	0.00
27 I	Chrysene-d12	0.400	0.400	0.0	99	0.00
28	Pyrene	0.400	0.392	2.0	98	0.00
29 S	Terphenyl-d14	0.400	0.397	0.8	98	0.00
30	Benzo(a)anthracene	0.400	0.408	-2.0	102	-0.01
31	Chrysene	0.400	0.410	-2.5	100	-0.01
32	Bis(2-ethylhexyl)phthalate	0.400	0.355	11.3	91	0.00
33	Indeno(1,2,3-cd)pyrene	0.400	0.406	-1.5	101	-0.03
34 I	Perylene-d12	0.400	0.400	0.0	99	-0.01
35	Benzo(b)fluoranthene	0.400	0.411	-2.7	102	-0.01
36	Benzo(k)fluoranthene	0.400	0.419	-4.7	103	-0.01
37 C	Benzo(a)pyrene	0.400	0.411	-2.7	102	-0.01
38	Dibenzo(a,h)anthracene	0.400	0.415	-3.7	102	-0.03
39	Benzo(g,h,i)perylene	0.400	0.416	-4.0	102	-0.03

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

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