

Data Path : Z:\HPCHEM1\BNA E\Data\BE041316\
 Data File : BE091657.D
 Acq On : 13 Apr 2016 14:59
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Apr 14 01:30:54 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE040816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 08 16:48:14 2016
 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	5.000	5.000	0.0	83	-0.02
2	1,4-Dioxane	0.400	0.397	0.8	84	0.00
3	n-Nitrosodimethylamine	0.400	0.354	11.5	79	0.00
4 S	2-Fluorophenol	0.400	0.355	11.3	78	0.00
5 S	Phenol-d6	0.400	0.346	13.5	79	0.00
6	bis(2-Chloroethyl)ether	0.400	0.375	6.3	83	0.00
7 I	Naphthalene-d8	5.000	5.000	0.0	85	-0.02
8 S	Nitrobenzene-d5	0.400	0.331	17.3	79	0.00
9	Nitrobenzene	0.400	0.429	-7.2	77	0.00
10	Naphthalene	0.400	0.400	0.0	86	0.00
11	Hexachlorobutadiene	0.400	0.394	1.5	84	0.00
12	2-Methylnaphthalene	0.400	0.389	2.8	85	-0.01
13 I	Acenaphthene-d10	5.000	5.000	0.0	83	0.00
14 S	2,4,6-Tribromophenol	0.400	0.476	-19.0	83	0.00
15 S	2-Fluorobiphenyl	0.400	0.405	-1.3	85	0.00
16	Acenaphthylene	0.400	0.375	6.3	96	0.00
17	Acenaphthene	0.400	0.414	-3.5	89	-0.01
18	Fluorene	0.400	0.390	2.5	84	-0.01
19 I	Phenanthrene-d10	5.000	5.000	0.0	90	-0.01
20	4-Bromophenyl-phenylether	0.400	0.345	13.8	81	0.00
21	Hexachlorobenzene	0.400	0.368	8.0	84	0.00
22	Pentachlorophenol	0.400	0.386	3.5	76	0.00
23	Phenanthrene	0.400	0.373	6.8	84	0.00
24	Anthracene	0.400	0.346	13.5	83	0.00
25	Fluoranthene	0.400	0.327	18.3	79	0.00
26 I	Chrysene-d12	5.000	5.000	0.0	77	0.00
27	Benzydine	0.400	0.483	-20.7	68	0.00
28	Pvrene	0.400	0.401	-0.3	83	-0.02
29 S	Terphenyl-d14	0.400	0.450	-12.5	88	0.00
30	Benzo(a)anthracene	0.400	0.411	-2.7	82	0.00
31	3,3'-Dichlorobenzidine	0.400	0.347	13.3	61	0.00
32	Chrysene	0.400	0.489	-22.2	96	-0.02
33	Bis(2-ethylhexyl)phthalate	0.400	0.345	13.8	42	0.00
34	Indeno(1,2,3-cd)pyrene	0.400	0.361	9.8	72	-0.02
35 I	Perylene-d12	5.000	5.000	0.0	79	-0.01
36	Benzo(b)fluoranthene	0.400	0.369	7.8	76	-0.01
37	Benzo(k)fluoranthene	0.400	0.394	1.5	81	-0.01
38 C	Benzo(a)pyrene	0.400	0.348	13.0	73	-0.01
39	Dibenzo(a,h)anthracene	0.400	0.350	12.5	72	-0.01
40	Benzo(a,h,i)perylene	0.400	0.364	9.0	74	-0.01

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

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