

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042616\
 Data File : BE091710.D
 Acq On : 26 Apr 2016 18:43
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Apr 27 01:04:29 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE042616.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 26 16:09:46 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	95	0.00
2	1,4-Dioxane	0.400	0.357	10.8	82	0.02
3	n-Nitrosodimethylamine	0.400	0.364	9.0	88	0.02
4 S	2-Fluorophenol	0.400	0.375	6.3	92	0.00
5 S	Phenol-d6	0.400	0.360	10.0	91	0.00
6	bis(2-Chloroethyl)ether	0.400	0.395	1.3	93	0.00
7 I	Naphthalene-d8	5.000	5.000	0.0	93	0.00
8 S	Nitrobenzene-d5	0.400	0.367	8.3	91	0.00
9	Nitrobenzene	0.400	0.359	10.3	90	0.00
10	Naphthalene	0.400	0.409	-2.2	94	0.00
11	Hexachlorobutadiene	0.400	0.408	-2.0	97	0.00
12	2-Methylnaphthalene	0.400	0.387	3.3	90	0.00
13 I	Acenaphthene-d10	5.000	5.000	0.0	92	0.00
14 S	2,4,6-Tribromophenol	0.400	0.444	-11.0	86	0.00
15 S	2-Fluorobiphenyl	0.400	0.394	1.5	91	0.00
16	Acenaphthylene	0.400	0.394	1.5	102	0.00
17	Acenaphthene	0.400	0.391	2.3	90	0.00
18	Fluorene	0.400	0.393	1.8	90	0.00
19 I	Phenanthrene-d10	5.000	5.000	0.0	90	0.00
20	4-Bromophenyl-phenylether	0.400	0.383	4.3	85	0.00
21	Hexachlorobenzene	0.400	0.395	1.3	89	0.00
22	Pentachlorophenol	0.400	0.360	10.0	81	0.00
23	Phenanthrene	0.400	0.411	-2.7	91	0.00
24	Anthracene	0.400	0.383	4.3	88	0.00
25	Fluoranthene	0.400	0.398	0.5	89	0.00
26 I	Chrysene-d12	5.000	5.000	0.0	81	0.00
27	Benzydine	0.400	0.423	-5.7	93	0.00
28	Pvrene	0.400	0.429	-7.2	90	0.00
29 S	Terphenyl-d14	0.400	0.430	-7.5	89	0.00
30	Benzo(a)anthracene	0.400	0.415	-3.7	86	0.00
31	3,3'-Dichlorobenzidine	0.400	0.394	1.5	92	-0.02
32	Chrysene	0.400	0.490	-22.5	94	0.00
33	Bis(2-ethylhexyl)phthalate	0.400	0.395	1.3	104	-0.02
34	Indeno(1,2,3-cd)pyrene	0.400	0.451	-12.7	90	0.00
35 I	Perylene-d12	5.000	5.000	0.0	91	0.00
36	Benzo(b)fluoranthene	0.400	0.404	-1.0	94	0.00
37	Benzo(k)fluoranthene	0.400	0.363	9.3	86	0.00
38 C	Benzo(a)pyrene	0.400	0.375	6.3	90	0.00
39	Dibenzo(a,h)anthracene	0.400	0.382	4.5	89	-0.01
40	Benzo(a,h,i)perylene	0.400	0.380	5.0	89	-0.01

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

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