

Data Path : Z:\HPCHEM1\BNA E\DATA\BE041816\
 Data File : BE091679.D
 Acq On : 19 Apr 2016 00:46
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Apr 19 05:47:16 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE041816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 18 17:20:40 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	91	0.00
2	1,4-Dioxane	0.400	0.378	5.5	94	0.00
3	n-Nitrosodimethylamine	0.400	0.420	-5.0	103	-0.02
4 S	2-Fluorophenol	0.400	0.434	-8.5	104	0.00
5 S	Phenol-d6	0.400	0.432	-8.0	107	0.00
6	bis(2-Chloroethyl)ether	0.400	0.406	-1.5	97	0.00
7 I	Naphthalene-d8	5.000	5.000	0.0	95	0.00
8 S	Nitrobenzene-d5	0.400	0.433	-8.2	119	0.00
9	Nitrobenzene	0.400	0.439	-9.7	118	0.00
10	Naphthalene	0.400	0.404	-1.0	95	0.00
11	Hexachlorobutadiene	0.400	0.419	-4.7	100	0.00
12	2-Methylnaphthalene	0.400	0.411	-2.7	100	0.00
13 I	Acenaphthene-d10	5.000	5.000	0.0	97	0.00
14 S	2,4,6-Tribromophenol	0.400	0.530	-32.5#	143	-0.01
15 S	2-Fluorobiphenyl	0.400	0.404	-1.0	99	0.00
16	Acenaphthylene	0.400	0.431	-7.7	120	0.00
17	Acenaphthene	0.400	0.421	-5.2	99	0.00
18	Fluorene	0.400	0.404	-1.0	100	0.00
19 I	Phenanthrene-d10	5.000	5.000	0.0	88	0.00
20	4-Bromophenyl-phenylether	0.400	0.405	-1.3	103	0.00
21	Hexachlorobenzene	0.400	0.403	-0.8	98	0.00
22	Pentachlorophenol	0.400	0.541	-35.3#	141	-0.02
23	Phenanthrene	0.400	0.408	-2.0	98	0.00
24	Anthracene	0.400	0.406	-1.5	104	-0.01
25	Fluoranthene	0.400	0.428	-7.0	109	-0.02
26 I	Chrysene-d12	5.000	5.000	0.0	92	0.00
27	Benzydine	0.400	0.469	-17.2	135	-0.02
28	Pvrene	0.400	0.516	-29.0#	126	0.00
29 S	Terphenyl-d14	0.400	0.487	-21.7	118	0.00
30	Benzo(a)anthracene	0.400	0.436	-9.0	106	0.00
31	3,3'-Dichlorobenzidine	0.400	0.585	-46.2#	165	-0.02
32	Chrysene	0.400	0.485	-21.2	110	0.00
33	Bis(2-ethylhexyl)phthalate	0.400	0.768	-92.0#	220	-0.02
34	Indeno(1,2,3-cd)pyrene	0.400	0.468	-17.0	112	-0.01
35 I	Perylene-d12	5.000	5.000	0.0	102	-0.01
36	Benzo(b)fluoranthene	0.400	0.442	-10.5	120	0.00
37	Benzo(k)fluoranthene	0.400	0.401	-0.3	107	-0.01
38 C	Benzo(a)pyrene	0.400	0.442	-10.5	122	-0.01
39	Dibenzo(a,h)anthracene	0.400	0.414	-3.5	112	-0.01
40	Benzo(a,h,i)perylene	0.400	0.409	-2.2	109	-0.01

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Compound	Amount	Calc.	%Dev Area	% Dev(min)
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(#) = Out of Range

SPCC's out = 0 CCC's out = 0