

Data Path : Z:\HPCHEM1\BNA E\DATA\BE070516\
 Data File : BE091982.D
 Acq On : 6 Jul 2016 4:06
 Operator : UM/SJ
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC0.4

Quant Time: Jul 06 06:52:07 2016
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE070516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 05 16:14:58 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	89	0.00
2	1,4-Dioxane	0.400	0.438	-9.5	93	0.00
3	n-Nitrosodimethylamine	0.400	0.367	8.3	90	0.00
4 S	2-Fluorophenol	0.400	0.376	6.0	89	0.00
5 S	Phenol-d6	0.400	0.350	12.5	86	0.00
6	bis(2-Chloroethyl)ether	0.400	0.392	2.0	94	0.00
7	n-Nitroso-di-n-propylamine	0.400	0.343	14.2	82	-0.02
8 I	Naphthalene-d8	5.000	5.000	0.0	89	0.00
9 S	Nitrobenzene-d5	0.400	0.366	8.5	83	-0.02
10	Nitrobenzene	0.400	0.360	10.0	86	0.00
11	bis(2-Chloroethoxy)methane	0.400	0.444	-11.0	86	0.00
12	Naphthalene	0.400	0.408	-2.0	89	0.00
13	Hexachlorobutadiene	0.400	0.404	-1.0	84	0.00
14	2-Methylnaphthalene	0.400	0.396	1.0	87	-0.01
15 I	Acenaphthene-d10	5.000	5.000	0.0	84	0.00
16 S	2,4,6-Tribromophenol	0.400	0.407	-1.7	84	0.00
17 S	2-Fluorobiphenyl	0.400	0.411	-2.7	84	0.00
18	Acenaphthylene	0.400	0.394	1.5	85	0.00
19	2,6-Dinitrotoluene	0.400	0.335	16.3	81	0.00
20	Acenaphthene	0.400	0.398	0.5	81	0.00
21	2,4-Dinitrotoluene	0.400	0.392	2.0	82	0.00
22	Fluorene	0.400	0.417	-4.2	85	0.00
23	4-Chlorophenyl-phenylether	0.400	0.414	-3.5	83	0.00
24 I	Phenanthrene-d10	5.000	5.000	0.0	89	0.00
25	4-Bromophenyl-phenylether	0.400	0.386	3.5	85	0.00
26	Hexachlorobenzene	0.400	0.424	-6.0	93	0.00
27	Pentachlorophenol	0.400	0.392	2.0	105	0.00
28	Phenanthrene	0.400	0.407	-1.7	88	0.00
29	Anthracene	0.400	0.388	3.0	88	0.00
30	Fluoranthene	0.400	0.402	-0.5	91	0.00
31 I	Chrysene-d12	5.000	5.000	0.0	102	0.00
32	Benzidine	0.400	0.379	5.3	60	0.00
33	Pyrene	0.400	0.366	8.5	93	0.00
34 S	Terphenyl-d14	0.400	0.357	10.8	90	0.00
35	Benzo(a)anthracene	0.400	0.456	-14.0	117	0.00
36	3,3'-Dichlorobenzidine	0.400	0.278	30.5#	73	0.00
37	Chrysene	0.400	0.464	-16.0	119	0.00
38	Bis(2-ethylhexyl)phthalate	0.400	0.314	21.5	87	0.00
39	Indeno(1,2,3-cd)pyrene	0.400	0.385	3.8	95	-0.02
40 I	Perylene-d12	5.000	5.000	0.0	101	0.00
41	Benzo(b)fluoranthene	0.400	0.399	0.3	101	0.00

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.382	4.5	98	0.00
43 C	Benzo(a)pyrene	0.400	0.381	4.8	97	0.00
44	Dibenzo(a,h)anthracene	0.400	0.374	6.5	96	0.00
45	Benzo(a,h,i)perylene	0.400	0.376	6.0	95	0.00

(#) = Out of Range

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