

Data Path : Z:\HPCHEM1\BNA E\DATA\BE093015\
 Data File : BE090985.D
 Acq On : 30 Sep 2015 18:58
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Oct 01 04:21:10 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE092815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 30 16:01:03 2015
 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	71	0.00
2	1,4-Dioxane	1.000	3.482	-248.2#	213	0.00
3	n-Nitrosodimethylamine	1.000	0.945	5.5	69	0.00
4 S	2-Fluorophenol	1.000	0.988	1.2	73	0.00
5 S	Phenol-d6	1.000	0.906	9.4	72	0.01
6	bis(2-Chloroethyl)ether	1.000	1.211	-21.1	78	0.00
7 I	Naphthalene-d8	5.000	5.000	0.0	75	0.00
8 S	Nitrobenzene-d5	1.000	0.887	11.3	71	0.00
9	Nitrobenzene	1.000	0.863	13.7	67	0.00
10	Naphthalene	1.000	0.994	0.6	74	0.00
11	Hexachlorobutadiene	1.000	0.930	7.0	68	0.00
12	2-Methylnaphthalene	1.000	0.941	5.9	75	0.00
13 I	Acenaphthene-d10	5.000	5.000	0.0	77	0.00
14 S	2,4,6-Tribromophenol	1.000	0.967	3.3	92	0.00
15 S	2-Fluorobiphenyl	1.000	0.862	13.8	72	0.00
16	Acenaphthylene	1.000	0.909	9.1	77	0.00
17	Acenaphthene	1.000	0.960	4.0	79	0.00
18	Fluorene	1.000	0.961	3.9	80	0.00
19 I	Phenanthrene-d10	5.000	5.000	0.0	83	0.00
20	4-Bromophenyl-phenylether	1.000	0.888	11.2	80	0.00
21	Hexachlorobenzene	1.000	0.971	2.9	80	0.00
22	Pentachlorophenol	1.000	0.899	10.1	91	0.00
23	Phenanthrene	1.000	0.966	3.4	84	-0.02
24	Anthracene	1.000	0.908	9.2	81	-0.02
25	Fluoranthene	1.000	0.972	2.8	88	0.00
26 I	Chrysene-d12	5.000	5.000	0.0	87	0.00
27	Benzydine	1.000	1.338	-33.8#	121	0.00
28	Pvrene	1.000	0.941	5.9	89	0.00
29 S	Terphenyl-d14	1.000	0.923	7.7	88	0.00
30	Benzo(a)anthracene	1.000	0.973	2.7	89	0.00
31	3,3'-Dichlorobenzidine	1.000	0.022	97.8#	0	-20.97#
32	Chrysene	1.000	0.962	3.8	86	0.00
33	Bis(2-ethylhexyl)phthalate	1.000	0.918	8.2	97	0.00
34	Indeno(1,2,3-cd)pyrene	1.000	1.003	-0.3	95	0.00
35 I	Perylene-d12	5.000	5.000	0.0	89	0.00
36	Benzo(b)fluoranthene	1.000	0.918	8.2	92	0.00
37	Benzo(k)fluoranthene	1.000	0.884	11.6	84	0.00
38 C	Benzo(a)pyrene	1.000	0.968	3.2	100	0.00
39	Dibenzo(a,h)anthracene	1.000	0.945	5.5	100	0.00
40	Benzo(a,h,i)perylene	1.000	0.951	4.9	98	0.00

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

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