

Data Path : Z:\HPCHEM1\BNA_E\Data\BE082515\
 Data File : BE090633.D
 Acq On : 26 Aug 2015 3:01
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDCCC001

Quant Time: Aug 26 04:31:39 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE081415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 14 16:52:42 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	99	0.00
2	1,4-Dioxane	1.000	1.202	-20.2	102	0.00
3	n-Nitrosodimethylamine	1.000	1.030	-3.0	104	0.00
4 S	2-Fluorophenol	1.000	0.739	26.1#	73	0.00
5 S	Phenol-d6	1.000	0.697	30.3#	72	0.00
6 I	Naphthalene-d8	5.000	5.000	0.0	91	0.00
7 S	Nitrobenzene-d5	1.000	0.914	8.6	87	0.00
8	Nitrobenzene	1.000	0.886	11.4	86	0.00
9	Naphthalene	1.000	0.971	2.9	89	0.00
10	Hexachlorobutadiene	1.000	1.044	-4.4	95	0.00
11	2-Methylnaphthalene	1.000	0.930	7.0	87	0.00
12 I	Acenaphthene-d10	5.000	5.000	0.0	99	0.00
13 S	2,4,6-Tribromophenol	1.000	0.830	17.0	91	0.00
14 S	2-Fluorobiphenyl	1.000	0.913	8.7	90	0.00
15	Acenaphthylene	1.000	0.931	6.9	94	0.00
16	Acenaphthene	1.000	0.980	2.0	98	0.00
17	Fluorene	1.000	1.055	-5.5	105	0.02
18 I	Phenanthrene-d10	5.000	5.000	0.0	116	0.00
19	4-Bromophenyl-phenylether	1.000	0.912	8.8	109	0.00
20	Hexachlorobenzene	1.000	0.976	2.4	114	-0.02
21	Pentachlorophenol	1.000	0.849	15.1	123	0.00
22	Phenanthrene	1.000	0.937	6.3	111	0.00
23	Anthracene	1.000	0.939	6.1	112	0.00
24	Fluoranthene	1.000	0.979	2.1	116	0.00
25 I	Chrysene-d12	5.000	5.000	0.0	133	0.00
26	Pyrene	1.000	0.888	11.2	121	0.00
27 S	Terphenyl-d14	1.000	0.877	12.3	118	0.00
28	Benzo(a)anthracene	1.000	0.938	6.2	131	0.00
29	Chrysene	1.000	0.997	0.3	134	0.00
30	Bis(2-ethylhexyl)phthalate	1.000	0.655	34.5#	93	0.00
31	Indeno(1,2,3-cd)pyrene	1.000	0.905	9.5	130	0.00
32 I	Perylene-d12	5.000	5.000	0.0	133	0.00
33	Benzo(b)fluoranthene	1.000	0.970	3.0	132	0.00
34	Benzo(k)fluoranthene	1.000	0.995	0.5	131	0.00
35 C	Benzo(a)pyrene	1.000	0.931	6.9	128	0.00
36	Dibenzo(a,h)anthracene	1.000	0.895	10.5	126	0.00
37	Benzo(g,h,i)perylene	1.000	0.913	8.7	124	0.00

(#) = Out of Range

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