

Data Path : Z:\HPCHEM1\BNA E\Data\BE021815\
 Data File : BE089634.D
 Acq On : 19 Feb 2015 12:40
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001

Quant Time: Feb 19 13:20:52 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SIMPAH-BE021815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 18 13:01:39 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	91	0.00
2	1,4-Dioxane	1.000	0.942	5.8	85	0.00
3	n-Nitrosodimethylamine	1.000	0.967	3.3	85	0.00
4 I	Naphthalene-d8	5.000	5.000	0.0	91	0.00
5 S	Nitrobenzene-d5	1.000	1.096	-9.6	103	0.00
6	Naphthalene	1.000	1.009	-0.9	91	0.00
7	2-Methylnaphthalene	1.000	1.023	-2.3	93	0.00
8 I	Acenaphthene-d10	5.000	5.000	0.0	97	0.00
9 S	2-Fluorobiphenyl	1.000	1.007	-0.7	95	0.00
10	Acenaphthylene	1.000	1.024	-2.4	97	0.00
11	Acenaphthene	1.000	0.997	0.3	95	0.00
12	Fluorene	1.000	1.016	-1.6	96	0.00
13 I	Phenanthrene-d10	5.000	5.000	0.0	99	0.00
14	Hexachlorobenzene	1.000	1.027	-2.7	101	0.00
15	Pentachlorophenol	1.000	0.943	5.7	92	0.00
16	Phenanthrene	1.000	0.995	0.5	98	0.00
17	Anthracene	1.000	1.032	-3.2	101	0.00
18	Fluoranthene	1.000	1.018	-1.8	101	0.00
19 I	Chrysene-d12	5.000	5.000	0.0	101	0.00
20	Pvrene	1.000	1.029	-2.9	101	0.00
21 S	Terphenyl-d14	1.000	1.028	-2.8	103	0.00
22	Benzo(a)anthracene	1.000	0.951	4.9	98	0.00
23	Chrysene	1.000	1.025	-2.5	103	0.00
24	Indeno(1,2,3-cd)pyrene	1.000	0.874	12.6	96	0.00
25 I	Perylene-d12	5.000	5.000	0.0	101	0.00
26	Benzo(b)fluoranthene	1.000	0.820	18.0	90	0.00
27	Benzo(k)fluoranthene	1.000	1.108	-10.8	106	0.00
28 C	Benzo(a)pyrene	1.000	0.847	15.3	93	0.00
29	Dibenzo(a,h)anthracene	1.000	0.869	13.1	95	0.00
30	Benzo(g,h,i)perylene	1.000	0.909	9.1	98	0.00

(#) = Out of Range

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