

Data Path : Z:\HPCHEM1\BNA_M\Data\BM050515\
 Data File : BM001215.D
 Acq On : 05 May 2015 19:28
 Operator : TP/IZ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM050515

Quant Time: May 06 13:07:18 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM050515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 06 12:46:35 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	101	0.00
2	1,4-Dioxane	1.000	1.068	-6.8	98	0.00
3	n-Nitrosodimethylamine	1.000	0.982	1.8	109	0.00
4 S	2-Fluorophenol	1.000	1.010	-1.0	99	0.00
5 S	Phenol-d6	1.000	1.001	-0.1	97	-0.02
6 I	Naphthalene-d8	5.000	5.000	0.0	100	0.00
7 S	Nitrobenzene-d5	1.000	0.996	0.4	96	0.00
8	Nitrobenzene	1.000	0.952	4.8	98	0.00
9	Naphthalene	1.000	1.030	-3.0	100	0.00
10	Hexachlorobutadiene	1.000	1.012	-1.2	101	0.00
11	2-Methylnaphthalene	1.000	1.019	-1.9	98	0.00
12 I	Acenaphthene-d10	5.000	5.000	0.0	98	0.00
13 S	2,4,6-Tribromophenol	1.000	0.923	7.7	98	0.00
14 S	2-Fluorobiphenyl	1.000	1.018	-1.8	98	-0.01
15	Acenaphthylene	1.000	1.007	-0.7	97	0.00
16	Acenaphthene	1.000	1.006	-0.6	97	0.00
17	Fluorene	1.000	1.046	-4.6	96	0.00
18 I	Phenanthrene-d10	5.000	5.000	0.0	97	0.00
19	Hexachlorobenzene	1.000	0.958	4.2	93	0.00
20	Pentachlorophenol	1.000	1.011	-1.1	93	0.00
21	Phenanthrene	1.000	1.036	-3.6	99	0.00
22	Anthracene	1.000	0.986	1.4	98	0.00
23	Fluoranthene	1.000	1.029	-2.9	96	0.00
24 I	Chrysene-d12	5.000	5.000	0.0	98	0.00
25	Pyrene	1.000	0.998	0.2	97	0.00
26 S	Terphenyl-d14	1.000	1.001	-0.1	97	0.00
27	Benzo(a)anthracene	1.000	1.011	-1.1	99	0.00
28	Chrysene	1.000	1.018	-1.8	97	0.00
29	Indeno(1,2,3-cd)pyrene	1.000	1.034	-3.4	99	-0.01
30 I	Perylene-d12	5.000	5.000	0.0	98	0.00
31	Benzo(b)fluoranthene	1.000	0.987	1.3	100	0.00
32	Benzo(k)fluoranthene	1.000	1.058	-5.8	97	0.00
33 C	Benzo(a)pyrene	1.000	1.015	-1.5	98	0.00
34	Dibenzo(a,h)anthracene	1.000	1.013	-1.3	98	0.00
35	Benzo(g,h,i)perylene	1.000	1.021	-2.1	99	0.00

(#) = Out of Range

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