

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030516\
 Data File : BM004566.D
 Acq On : 05 Mar 2016 21:08
 Operator : SJ/UM
 Sample : SSTDICV001
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM030516

Quant Time: Mar 07 07:58:25 2016
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BM030516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 07 07:54:56 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	119	0.00
2	1,4-Dioxane	1.000	0.967	3.3	120	0.00
3	n-Nitrosodimethylamine	1.000	0.906	9.4	120	0.00
4 S	2-Fluorophenol	1.000	0.896	10.4	119	0.00
5 S	Phenol-d6	1.000	0.917	8.3	121	0.00
6	bis(2-Chloroethyl)ether	1.000	0.957	4.3	123	0.00
7 I	Naphthalene-d8	5.000	5.000	0.0	117	0.00
8 S	Nitrobenzene-d5	1.000	0.926	7.4	121	0.00
9	Nitrobenzene	1.000	0.907	9.3	121	0.00
10	Naphthalene	1.000	1.006	-0.6	132	0.00
11	Hexachlorobutadiene	1.000	0.941	5.9	122	0.00
12	2-Methylnaphthalene	1.000	0.936	6.4	123	0.00
13 I	Acenaphthene-d10	5.000	5.000	0.0	116	0.00
14 S	2,4,6-Tribromophenol	1.000	0.923	7.7	120	0.00
15 S	2-Fluorobiphenyl	1.000	0.995	0.5	123	0.00
16	Acenaphthylene	1.000	0.963	3.7	121	0.00
17	Acenaphthene	1.000	1.009	-0.9	125	0.00
18	Fluorene	1.000	0.976	2.4	122	0.00
19 I	Phenanthrene-d10	5.000	5.000	0.0	120	0.00
20	Hexachlorobenzene	1.000	0.965	3.5	121	0.00
21	Pentachlorophenol	1.000	0.741	25.9#	113	0.00
22	Phenanthrene	1.000	1.006	-0.6	127	0.00
23	Anthracene	1.000	0.931	6.9	126	0.00
24	Fluoranthene	1.000	0.959	4.1	123	0.00
25 I	Chrysene-d12	5.000	5.000	0.0	117	0.00
26	Benzidine	1.000	1.252	-25.2#	213	-0.02
27	Pyrene	1.000	0.981	1.9	125	0.00
28 S	Terphenyl-d14	1.000	0.962	3.8	125	0.00
29	Benzo(a)anthracene	1.000	0.951	4.9	123	0.00
30	3,3'-Dichlorobenzidine	1.000	0.992	0.8	126	0.00
31	Chrysene	1.000	1.020	-2.0	126	-0.02
32	Bis(2-ethylhexyl)phthalate	1.000	0.847	15.3	113	0.00
33	Indeno(1,2,3-cd)pyrene	1.000	0.969	3.1	124	0.00
34 I	Perylene-d12	5.000	5.000	0.0	122	0.00
35 C	Benzo(a)pyrene	1.000	0.939	6.1	126	-0.01
36	Dibenzo(a,h)anthracene	1.000	0.899	10.1	125	0.00
37	Benzo(g,h,i)perylene	1.000	0.900	10.0	125	0.00

(#) = Out of Range

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