

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101515\
 Data File : BM002915.D
 Acq On : 15 Oct 2015 15:26
 Operator : UM/IZ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ICVBM101515

Quant Time: Oct 16 06:51:15 2015
 Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM101515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 15 16:21:18 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	78	0.00
2	1,4-Dioxane	1.000	0.959	4.1	77	0.00
3	n-Nitrosodimethylamine	1.000	1.022	-2.2	78	-0.02
4 S	2-Fluorophenol	1.000	0.947	5.3	76	-0.02
5 S	Phenol-d6	1.000	0.998	0.2	77	-0.02
6	bis(2-Chloroethyl)ether	1.000	1.002	-0.2	77	-0.02
7 I	Naphthalene-d8	5.000	5.000	0.0	80	0.00
8 S	Nitrobenzene-d5	1.000	0.965	3.5	77	-0.02
9	Nitrobenzene	1.000	0.959	4.1	76	-0.02
10	Naphthalene	1.000	0.968	3.2	79	0.00
11	Hexachlorobutadiene	1.000	0.961	3.9	78	-0.01
12	2-Methylnaphthalene	1.000	0.990	1.0	80	0.00
13 I	Acenaphthene-d10	5.000	5.000	0.0	80	0.00
14 S	2,4,6-Tribromophenol	1.000	0.924	7.6	80	0.00
15 S	2-Fluorobiphenyl	1.000	0.995	0.5	81	0.00
16	Acenaphthylene	1.000	0.985	1.5	80	0.00
17	Acenaphthene	1.000	0.968	3.2	81	0.00
18	Fluorene	1.000	0.993	0.7	81	0.00
19 I	Phenanthrene-d10	5.000	5.000	0.0	90	0.00
20	4-Bromophenyl-phenylether	1.000	1.078	-7.8	100	0.00
21	Hexachlorobenzene	1.000	1.094	-9.4	79	0.02
22	Pentachlorophenol	1.000	0.979	2.1	89	0.00
23	Phenanthrene	1.000	0.935	6.5	85	0.02
24	Anthracene	1.000	0.850	15.0	76	0.00
25	Fluoranthene	1.000	0.974	2.6	90	0.00
26 I	Chrysene-d12	5.000	5.000	0.0	87	0.01
27	Benzydine	1.000	0.944	5.6	96	0.00
28	Pvrene	1.000	0.995	0.5	88	0.00
29 S	Terphenyl-d14	1.000	0.993	0.7	88	0.00
30	Benzo(a)anthracene	1.000	0.915	8.5	77	0.01
31	3,3'-Dichlorobenzidine	1.000	0.936	6.4	86	0.01
32	Chrysene	1.000	0.994	0.6	95	0.00
33	Bis(2-ethylhexyl)phthalate	1.000	0.890	11.0	70	0.00
34	Indeno(1,2,3-cd)pyrene	1.000	1.011	-1.1	96	0.00
35 I	Perylene-d12	5.000	5.000	0.0	97	0.00
36	Benzo(b)fluoranthene	1.000	1.011	-1.1	97	0.00
37	Benzo(k)fluoranthene	1.000	0.935	6.5	93	0.00
38 C	Benzo(a)pyrene	1.000	0.952	4.8	94	0.00
39	Dibenzo(a,h)anthracene	1.000	0.960	4.0	97	0.00
40	Benzo(a,h,i)perylene	1.000	0.978	2.2	99	0.00

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

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