Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM040715\
Data File : BM000900.D
Acq On : 08 Apr 2015 04:20
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDCCC001EC

Quant Time: Apr 08 06:04:23 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-BM040715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 07 15:59:39 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	109	-0.04
2		1,4-Dioxane	1.000	1.013	-1.3	104	0.00
3		n-Nitrosodimethylamine	1.000	0.939	6.1	111	0.00
4	S	2-Fluorophenol	1.000	0.986	1.4	111	0.00
5	S	Phenol-d6	1.000	0.937	6.3	113	0.00
6		2-Chlorophenol	1.000	0.986	1.4	122	0.00
7	C	Phenol	1.000	0.858	14.2	99	0.00
8		bis(2-Chloroethyl)ether	1.000	0.984	1.6	125	0.00
9		2-Methylphenol	1.000	0.851	14.9	95	0.00
10		3+4-Methylphenols	1.000	0.905	9.5	108	0.00
11	I	Naphthalene-d8	5.000	5.000	0.0	108	-0.03
12	S	Nitrobenzene-d5	1.000	0.966	3.4	112	-0.04
13		Nitrobenzene	1.000	0.886	11.4	99	0.00
14		2-Nitrophenol	1.000	0.791	20.9	84	0.00
15		2,4-Dimethylphenol	1.000	0.924	7.6	101	0.00
16	C	2,4-Dichlorophenol	1.000	0.967	3.3	116	0.00
17	C	Hexachlorobutadiene	1.000	1.095	-9.5	117	0.00
18		4-Chloro-3-methylphenol	1.000	0.953	4.7	112	0.00
19	I	Acenaphthene-d10	5.000	5.000	0.0	109	0.00
20	S	2,4,6-Tribromophenol	1.000	0.900	10.0	103	0.02
21		2,4,6-Trichlorophenol	1.000	0.946	5.4	134	0.00
22		2,4,5-Trichlorophenol	1.000	0.953	4.7	127	0.00
23	S	2-Fluorobiphenyl	1.000	0.996	0.4	109	0.00
24	P	2,4-Dinitrophenol	1.000	0.000	100.0#	0	-14.40#
25		4-Nitrophenol	1.000	0.882	11.8	95	-0.01
26	I	Phenanthrene-d10	5.000	5.000	0.0	124	0.00
27		4,6-Dinitro-2-methylphenol	1.000	0.773	22.7	94	0.00
28		Hexachlorobenzene	1.000	0.855	14.5	85	0.00
29	C	Pentachlorophenol	1.000	0.964	3.6	166	0.00
30	I	Chrysene-d12	5.000	5.000	0.0	106	0.00
31		Benzidine	1.000	0.942	5.8	102	0.00
32	S	Terphenyl-d14	1.000	1.002	-0.2	106	0.00
33		3,3'-Dichlorobenzidine	1.000	0.929	7.1	100	0.00
34	I	Perylene-d12	5.000	5.000	0.0	105	0.00

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