Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM120715\
Data File : BM003407.D
Acq On : 07 Dec 2015 14:48
Operator : UM/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM120715

Quant Time: Dec 07 15:30:40 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-BM120715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 07 15:08:30 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	114	0.00
2		1,4-Dioxane	1.000	0.952	4.8	109	0.00
3		n-Nitrosodimethylamine	1.000	0.978	2.2	114	0.00
4	S	2-Fluorophenol	1.000	0.970	3.0	118	-0.02
5	S	Phenol-d6	1.000	1.001	-0.1	120	0.00
6		2-Chlorophenol	1.000	1.006	-0.6	119	0.00
7	C	Phenol	1.000	0.986	1.4	118	0.00
8		bis(2-Chloroethyl)ether	1.000	1.007	-0.7	118	0.00
9		2-Methylphenol	1.000	1.003	-0.3	121	0.00
10		3+4-Methylphenols	1.000	1.005	-0.5	119	0.00
11	I	Naphthalene-d8	5.000	5.000	0.0	116	0.00
12	S	Nitrobenzene-d5	1.000	1.024	-2.4	122	0.00
13		Nitrobenzene	1.000	1.014	-1.4	123	0.00
14	C	2-Nitrophenol	1.000	0.981	1.9	124	0.00
15		2,4-Dimethylphenol	1.000	1.019	-1.9	122	0.00
16	C	2,4-Dichlorophenol	1.000	1.020	-2.0	119	0.00
17		Hexachlorobutadiene	1.000	0.955	4.5	113	0.00
18	C	4-Chloro-3-methylphenol	1.000	0.984	1.6	119	0.03
19		2-Methylnaphthalene	1.000	0.989	1.1	118	0.00
20	I	Acenaphthene-d10	5.000	5.000	0.0	145	0.00
21	S	2,4,6-Tribromophenol	1.000	0.765	23.5	101	0.00
22	C	2,4,6-Trichlorophenol	1.000	0.835	16.5	127	0.00
23		2,4,5-Trichlorophenol	1.000	0.813	18.7	116	0.00
24	S	2-Fluorobiphenyl	1.000	0.872	12.8	124	0.00
25	P	2,4-Dinitrophenol	1.000	1.038	-3.8	137	0.00
26	P	4-Nitrophenol	1.000	0.894	10.6	120	0.00
27	I	Phenanthrene-d10	5.000	5.000	0.0	133	0.00
28		4,6-Dinitro-2-methylphenol	1.000	1.058	-5.8	147	0.00
29		Hexachlorobenzene	1.000	0.858	14.2	107	0.00
30	C	Pentachlorophenol	1.000	0.854	14.6	120	0.00
31	I	Chrysene-d12	5.000	5.000	0.0	85	0.00
32		Benzidine	1.000	1.145	-14.5	126	0.00
33	S	Terphenyl-d14	1.000	1.175	-17.5	104	0.00
34		3,3'-Dichlorobenzidine	1.000	0.971	2.9	134	0.00
35	I	Perylene-d12	5.000	5.000	0.0	114	0.01

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