Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101315\
Data File : BM002899.D
Acq On : 13 Oct 2015 22:48
Operator : UM/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM101315

Quant Time: Oct 14 17:42:48 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-ACID-BM0101315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 14 17:36:42 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	87	0.00
2	S	2-Fluorophenol	1.000	0.855	14.5	78	0.02
3	S	Phenol-d6	1.000	0.874	12.6	80	0.02
4		2-Chlorophenol	1.000	0.903	9.7	80	0.00
5	C	Phenol	1.000	0.911	8.9	80	0.00
6		2-Methylphenol	1.000	0.888	11.2	80	0.00
7		3+4-Methylphenols	1.000	0.887	11.3	78	0.01
8	I	Naphthalene-d8	5.000	5.000	0.0	85	0.00
9	S	Nitrobenzene-d5	1.000	0.944	5.6	81	0.00
10	C	2-Nitrophenol	1.000	0.864	13.6	77	0.00
11		2,4-Dimethylphenol	1.000	0.904	9.6	79	0.00
12	C	2,4-Dichlorophenol	1.000	0.919	8.1	76	0.02
13	C	4-Chloro-3-methylphenol	1.000	0.882	11.8	78	0.00
14	I	Acenaphthene-d10	5.000	5.000	0.0	86	0.00
15	S	2,4,6-Tribromophenol	1.000	0.860	14.0	77	0.00
16	C	2,4,6-Trichlorophenol	1.000	0.849	15.1	75	0.00
17		2,4,5-Trichlorophenol	1.000	0.878	12.2	77	0.00
18	S	2-Fluorobiphenyl	1.000	0.926	7.4	81	0.01
19	P	2,4-Dinitrophenol	1.000	0.818	18.2	55	0.02
20	P	4-Nitrophenol	1.000	0.877	12.3	69	0.00
21	I	Phenanthrene-d10	5.000	5.000	0.0	90	0.00
22		4,6-Dinitro-2-methylphenol	1.000	0.869	13.1	73	0.00
23	C	Pentachlorophenol	1.000	0.965	3.5	69	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	93	0.00
25	S	Terphenyl-d14	1.000	0.898	10.2	86	0.00
26	I	Perylene-d12	5.000	5.000	0.0	91	0.00

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