Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101415\
Data File : BM002901.D
Acq On : 14 Oct 2015 16:38
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Oct 15 01:02:13 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	92	-0.01
2	S	2-Fluorophenol	1.000	0.898	10.2	87	0.02
3	S	Phenol-d6	1.000	0.856	14.4	83	0.02
4		2-Chlorophenol	1.000	0.889	11.1	83	0.00
5	C	Phenol	1.000	0.892	10.8	83	0.00
6		2-Methylphenol	1.000	0.861	13.9	82	0.00
7		3+4-Methylphenols	1.000	0.871	12.9	81	0.00
8	I	Naphthalene-d8	5.000	5.000	0.0	90	0.00
9	S	Nitrobenzene-d5	1.000	0.920	8.0	83	0.00
10	C	2-Nitrophenol	1.000	0.838	16.2	78	0.00
11		2,4-Dimethylphenol	1.000	0.881	11.9	81	0.00
12	C	2,4-Dichlorophenol	1.000	0.885	11.5	77	0.02
13	C	4-Chloro-3-methylphenol	1.000	0.848	15.2	79	0.00
14	I	Acenaphthene-d10	5.000	5.000	0.0	89	0.00
15	S	2,4,6-Tribromophenol	1.000	0.842	15.8	78	0.00
16	C	2,4,6-Trichlorophenol	1.000	0.824	17.6	75	0.00
17		2,4,5-Trichlorophenol	1.000	0.856	14.4	77	0.00
18	S	2-Fluorobiphenyl	1.000	0.930	7.0	83	0.00
19	P	2,4-Dinitrophenol	1.000	0.836	16.4	60	0.02
20	P	4-Nitrophenol	1.000	0.847	15.3	66	0.00
21	I	Phenanthrene-d10	5.000	5.000	0.0	93	0.00
22		4,6-Dinitro-2-methylphenol	1.000	0.874	12.6	76	0.00
23	C	Pentachlorophenol	1.000	0.960	4.0	71	0.00
24	I	Chrysene-d12	5.000	5.000	0.0	94	0.00
25	S	Terphenyl-d14	1.000	0.902	9.8	87	0.00
26	I	Perylene-d12	5.000	5.000	0.0	93	0.00

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