Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM040315\
Data File : BM000819.D
Acq On : 02 Apr 2015 22:44
Operator : TP/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM040315

Quant Time: Apr 03 18:28:18 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SIM-BM040315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 03 18:16:53 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	116	0.00
2		1,4-Dioxane	1.000	0.979	2.1	111	0.00
3		n-Nitrosodimethylamine	1.000	1.045	-4.5	121	-0.02
4	S	2-Fluorophenol	1.000	1.042	-4.2	120	0.00
5	S	Phenol-d6	1.000	0.998	0.2	122	0.00
6		2-Chlorophenol	1.000	1.004	-0.4	121	0.00
7	C	Phenol	1.000	1.034	-3.4	123	0.00
8		bis(2-Chloroethyl)ether	1.000	0.978	2.2	122	0.00
9		2-Methylphenol	1.000	1.021	-2.1	123	0.00
10		3+4-Methylphenols	1.000	0.970	3.0	118	0.00
11	I	Naphthalene-d8	5.000	5.000	0.0	118	0.00
12	S	Nitrobenzene-d5	1.000	0.993	0.7	122	0.00
13		Nitrobenzene	1.000	1.040	-4.0	126	0.00
14		2-Nitrophenol	1.000	1.030	-3.0	122	0.00
15		2,4-Dimethylphenol	1.000	1.052	-5.2	122	0.00
16	C	2,4-Dichlorophenol	1.000	1.001	-0.1	121	0.00
17	C	Hexachlorobutadiene	1.000	1.023	-2.3	118	0.00
18		4-Chloro-3-methylphenol	1.000	1.005	-0.5	120	0.00
19	I	Acenaphthene-d10	5.000	5.000	0.0	118	0.00
20	S	2,4,6-Tribromophenol	1.000	0.948	5.2	119	0.00
21		2,4,6-Trichlorophenol	1.000	0.971	2.9	116	0.00
22		2,4,5-Trichlorophenol	1.000	1.070	-7.0	119	0.00
23	S	2-Fluorobiphenyl	1.000	1.043	-4.3	120	0.00
24	P	2,4-Dinitrophenol	1.000	0.452	54.8#	115	0.00
25		4-Nitrophenol	1.000	1.041	-4.1	129	0.00
26	I	Phenanthrene-d10	5.000	5.000	0.0	111	0.00
27		4,6-Dinitro-2-methylphenol	1.000	0.892	10.8	102	0.03
28		Hexachlorobenzene	1.000	1.046	-4.6	117	0.00
29	C	Pentachlorophenol	1.000	1.031	-3.1	114	0.00
30	I	Chrysene-d12	5.000	5.000	0.0	116	0.00
31		Benzidine	1.000	1.182	-18.2	777	-0.02
32	S	Terphenyl-d14	1.000	1.041	-4.1	118	0.00
33		3,3'-Dichlorobenzidine	1.000	0.959	4.1	148	0.00
34	I	Perylene-d12	5.000	5.000	0.0	115	0.00

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