Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM040715\
Data File : BM000881.D
Acq On : 07 Apr 2015 15:50
Operator : TP/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM040715

Quant Time: Apr 07 16:37:48 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	145	0.00
2		1,4-Dioxane	1.000	1.040	-4.0	142	0.00
3		n-Nitrosodimethylamine	1.000	1.035	-3.5	167	-0.02
4	S	2-Fluorophenol	1.000	1.055	-5.5	158	0.00
5	S	Phenol-d6	1.000	1.027	-2.7	168	0.00
6		2-Chlorophenol	1.000	1.001	-0.1	165	0.00
7	C	Phenol	1.000	1.039	-3.9	159	0.00
8		bis(2-Chloroethyl)ether	1.000	1.002	-0.2	170	0.00
9		2-Methylphenol	1.000	1.048	-4.8	155	0.00
10		3+4-Methylphenols	1.000	1.012	-1.2	166	0.00
11	I	Naphthalene-d8	5.000	5.000	0.0	150	0.00
12	S	Nitrobenzene-d5	1.000	1.011	-1.1	163	-0.04
13		Nitrobenzene	1.000	1.027	-2.7	159	0.00
14		2-Nitrophenol	1.000	1.008	-0.8	155	0.00
15		2,4-Dimethylphenol	1.000	1.037	-3.7	158	0.00
16	C	2,4-Dichlorophenol	1.000	1.000	0.0	168	0.00
17	C	Hexachlorobutadiene	1.000	1.038	-3.8	155	0.00
18		4-Chloro-3-methylphenol	1.000	0.997	0.3	164	0.00
19	I	Acenaphthene-d10	5.000	5.000	0.0	148	0.00
20	S	2,4,6-Tribromophenol	1.000	0.965	3.5	155	0.02
21		2,4,6-Trichlorophenol	1.000	0.959	4.1	185	0.00
22		2,4,5-Trichlorophenol	1.000	0.980	2.0	179	0.00
23	S	2-Fluorobiphenyl	1.000	1.054	-5.4	156	0.00
24	P	2,4-Dinitrophenol	1.000	0.000	100.0#	0	-14.40#
25		4-Nitrophenol	1.000	1.169	-16.9	196	-0.01
26	I	Phenanthrene-d10	5.000	5.000	0.0	165	0.00
27		4,6-Dinitro-2-methylphenol	1.000	0.989	1.1	174	0.00
28		Hexachlorobenzene	1.000	0.937	6.3	123	0.02
29	C	Pentachlorophenol	1.000	1.087	-8.7	284	0.00
30	I	Chrysene-d12	5.000	5.000	0.0	143	0.00
31		Benzidine	1.000	1.201	-20.1	189	0.00
32	S	Terphenyl-d14	1.000	1.043	-4.3	150	0.00
33		3,3'-Dichlorobenzidine	1.000	1.155	-15.5	179	0.00
34	I	Perylene-d12	5.000	5.000	0.0	146	0.00

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