Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM040315\
Data File : BM000839.D
Acq On : 03 Apr 2015 13:06
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDCCC001

Quant Time: Apr 03 18:52:16 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	99	0.00
2		1,4-Dioxane	0.756	0.723	4.4	103	0.00
3		n-Nitrosodimethylamine	0.635	0.648	-2.0	100	0.00
4	S	2-Fluorophenol	1.377	1.448	-5.2	103	0.00
5	S	Phenol-d6	1.921	2.065	-7.5	105	0.00
6		2-Chlorophenol	1.365	1.481	-8.5	107	0.00
7	C	Phenol	2.063	2.080	-0.8	101	0.00
8		bis(2-Chloroethyl)ether	1.435	1.484	-3.4	108	0.00
9		2-Methylphenol	1.472	1.516	-3.0	101	0.00
10		3+4-Methylphenols	1.558	1.726	-10.8	100	0.00
11	I	Naphthalene-d8	1.000	1.000	0.0	101	0.00
12	S	Nitrobenzene-d5	0.233	0.226	3.0	102	0.00
13		Nitrobenzene	0.272	0.255	6.3	98	0.00
14		2-Nitrophenol	0.142	0.129	9.2	98	0.00
15		2,4-Dimethylphenol	0.283	0.289	-2.1	101	0.00
16	C	2,4-Dichlorophenol	0.251	0.257	-2.4	101	0.00
17	C	Hexachlorobutadiene	0.193	0.203	-5.2	104	-0.03
18		4-Chloro-3-methylphenol	0.251	0.253	-0.8	103	0.00
19	I	Acenaphthene-d10	1.000	1.000	0.0	105	0.00
20	S	2,4,6-Tribromophenol	0.149	0.148	0.7	94	0.00
21		2,4,6-Trichlorophenol	0.270	0.277	-2.6	89	0.00
22		2,4,5-Trichlorophenol	0.331	0.363	-9.7	96	0.00
23	S	2-Fluorobiphenyl	1.540	1.618	-5.1	107	0.00
24	P	2,4-Dinitrophenol	0.049	0.010#	79.6#	48#	0.00
25		4-Nitrophenol	0.098	0.075	23.5	82	0.00
26	I	Phenanthrene-d10	1.000	1.000	0.0	109	0.00
27		4,6-Dinitro-2-methylphenol	0.044	0.035	20.5	81	0.00
28		Hexachlorobenzene	0.282	0.267	5.3	101	0.00
29	C	Pentachlorophenol	0.049	0.028	42.9#	58	0.00
30	I	Chrysene-d12	1.000	1.000	0.0	102	0.00
31		Benzidine	0.201	0.223	-10.9	898#	-0.02
32	S	Terphenyl-d14	0.602	0.624	-3.7	104	0.00
33		3,3'-Dichlorobenzidine	0.336	0.337	-0.3	123	0.00
34	I	Perylene-d12	1.000	1.000	0.0	103	0.00

(#) = Out of Range SPCC's out = 1 CCC's out = 1