Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM040715\
Data File : BM000887.D
Acq On : 07 Apr 2015 19:25
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDCCC001EC

Quant Time: Apr 08 01:55:52 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	108	0.00
2		1,4-Dioxane	0.864	0.834	3.5	107	0.00
3		n-Nitrosodimethylamine	0.746	0.754	-1.1	116	-0.02
4	S	2-Fluorophenol	1.678	1.681	-0.2	112	0.00
5	S	Phenol-d6	2.645	2.663	-0.7	116	0.00
6		2-Chlorophenol	2.076	2.122	-2.2	118	0.00
7	C	Phenol	2.091	1.858	11.1	102	0.00
8		bis(2-Chloroethyl)ether	2.038	2.132	-4.6	122	0.00
9		2-Methylphenol	1.533	1.460	4.8	105	0.00
10		3+4-Methylphenols	1.999	1.951	2.4	113	0.00
11	I	Naphthalene-d8	1.000	1.000	0.0	112	0.00
12	S	Nitrobenzene-d5	0.249	0.234	6.0	113	-0.04
13		Nitrobenzene	0.251	0.229	8.8	105	0.00
14		2-Nitrophenol	0.158	0.140	11.4	113	0.00
15		2,4-Dimethylphenol	0.287	0.278	3.1	110	0.00
16	C	2,4-Dichlorophenol	0.281	0.274	2.5	115	0.00
17	C	Hexachlorobutadiene	0.228	0.232	-1.8	113	0.00
18		4-Chloro-3-methylphenol	0.274	0.264	3.6	115	0.00
19	I	Acenaphthene-d10	1.000	1.000	0.0	113	0.00
20	S	2,4,6-Tribromophenol	0.145	0.126	13.1	102	0.02
21		2,4,6-Trichlorophenol	0.256	0.266	-3.9	124	0.00
22		2,4,5-Trichlorophenol	0.350	0.348	0.6	123	0.00
23	S	2-Fluorobiphenyl	1.541	1.543	-0.1	114	0.00
24	P	2,4-Dinitrophenol	0.000	0.000#	0.0	0#	-14.40#
25		4-Nitrophenol	0.061	0.070	-14.8	107	-0.01
26	I	Phenanthrene-d10	1.000	1.000	0.0	125	0.00
27		4,6-Dinitro-2-methylphenol	0.049	0.035	28.6#	96	0.00
28		Hexachlorobenzene	0.222	0.197	11.3	87	0.02
29	C	Pentachlorophenol	0.052	0.025	51.9#	145	0.00
30	I	Chrysene-d12	1.000	1.000	0.0	104	0.00
31		Benzidine	0.222	0.229	-3.2	108	0.00
32	S	Terphenyl-d14	0.617	0.618	-0.2	104	0.00
33		3,3'-Dichlorobenzidine	0.330	0.348	-5.5	110	0.00
34	I	Perylene-d12	1.000	1.000	0.0	104	0.00

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