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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	145	0.00
2		1,4-Dioxane	0.864	0.829	4.1	142	0.00
3		n-Nitrosodimethylamine	0.746	0.814	-9.1	167#	-0.02
4	S	2-Fluorophenol	1.678	1.771	-5.5	158#	0.00
5	S	Phenol-d6	2.645	2.891	-9.3	168#	0.00
6		2-Chlorophenol	2.076	2.223	-7.1	165#	0.00
7	C	Phenol	2.091	2.173	-3.9	159#	0.00
8		bis(2-Chloroethyl)ether	2.038	2.228	-9.3	170#	0.00
9		2-Methylphenol	1.533	1.607	-4.8	155#	0.00
10		3+4-Methylphenols	1.999	2.142	-7.2	166#	0.00
11	I	Naphthalene-d8	1.000	1.000	0.0	150	0.00
12	S	Nitrobenzene-d5	0.249	0.252	-1.2	163#	-0.04
13		Nitrobenzene	0.251	0.258	-2.8	159#	0.00
14		2-Nitrophenol	0.158	0.143	9.5	155#	0.00
15		2,4-Dimethylphenol	0.287	0.297	-3.5	158#	0.00
16	C	2,4-Dichlorophenol	0.281	0.300	-6.8	168#	0.00
17	C	Hexachlorobutadiene	0.228	0.237	-3.9	155#	0.00
18		4-Chloro-3-methylphenol	0.274	0.280	-2.2	164#	0.00
19	I	Acenaphthene-d10	1.000	1.000	0.0	148	0.00
20	S	2,4,6-Tribromophenol	0.145	0.146	-0.7	155#	0.02
21		2,4,6-Trichlorophenol	0.256	0.302	-18.0	185#	0.00
22		2,4,5-Trichlorophenol	0.350	0.387	-10.6	179#	0.00
23	S	2-Fluorobiphenyl	1.541	1.625	-5.5	156#	0.00
24	P	2,4-Dinitrophenol	0.000	0.000#	0.0	0#	-14.40#
25		4-Nitrophenol	0.061	0.097	-59.0#	196#	-0.01
26	I	Phenanthrene-d10	1.000	1.000	0.0	165#	0.00
27		4,6-Dinitro-2-methylphenol	0.049	0.047	4.1	174#	0.00
28		Hexachlorobenzene	0.222	0.209	5.9	123	0.02
29	C	Pentachlorophenol	0.052	0.037	28.8#	284#	0.00
30	I	Chrysene-d12	1.000	1.000	0.0	143	0.00
31		Benzidine	0.222	0.291	-31.1#	189#	0.00
32	S	Terphenyl-d14	0.617	0.643	-4.2	150	0.00
33		3,3'-Dichlorobenzidine	0.330	0.410	-24.2	179#	0.00
34	I	Perylene-d12	1.000	1.000	0.0	146	0.00

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