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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	116	0.00
2		1,4-Dioxane	0.756	0.665	12.0	111	0.00
3		n-Nitrosodimethylamine	0.635	0.664	-4.6	121	-0.02
4	S	2-Fluorophenol	1.377	1.435	-4.2	120	0.00
5	S	Phenol-d6	1.921	2.040	-6.2	122	0.00
6		2-Chlorophenol	1.365	1.420	-4.0	121	0.00
7	C	Phenol	2.063	2.132	-3.3	123	0.00
8		bis(2-Chloroethyl)ether	1.435	1.420	1.0	122	0.00
9		2-Methylphenol	1.472	1.562	-6.1	123	0.00
10		3+4-Methylphenols	1.558	1.715	-10.1	118	0.00
11	I	Naphthalene-d8	1.000	1.000	0.0	118	0.00
12	S	Nitrobenzene-d5	0.233	0.232	0.4	122	0.00
13		Nitrobenzene	0.272	0.283	-4.0	126	0.00
14		2-Nitrophenol	0.142	0.137	3.5	122	0.00
15		2,4-Dimethylphenol	0.283	0.298	-5.3	122	0.00
16	C	2,4-Dichlorophenol	0.251	0.263	-4.8	121	0.00
17	C	Hexachlorobutadiene	0.193	0.198	-2.6	118	0.00
18		4-Chloro-3-methylphenol	0.251	0.252	-0.4	120	0.00
19	I	Acenaphthene-d10	1.000	1.000	0.0	118	0.00
20	S	2,4,6-Tribromophenol	0.149	0.166	-11.4	119	0.00
21		2,4,6-Trichlorophenol	0.270	0.322	-19.3	116	0.00
22		2,4,5-Trichlorophenol	0.331	0.397	-19.9	119	0.00
23	S	2-Fluorobiphenyl	1.540	1.606	-4.3	120	0.00
24	P	2,4-Dinitrophenol	0.049	0.022#	55.1#	115	0.00
25		4-Nitrophenol	0.098	0.104	-6.1	129	0.00
26	I	Phenanthrene-d10	1.000	1.000	0.0	111	0.00
27		4,6-Dinitro-2-methylphenol	0.044	0.043	2.3	102	0.03
28		Hexachlorobenzene	0.282	0.304	-7.8	117	0.00
29	C	Pentachlorophenol	0.049	0.054	-10.2	114	0.00
30	I	Chrysene-d12	1.000	1.000	0.0	116	0.00
31		Benzidine	0.201	0.171	14.9	777#	-0.02
32	S	Terphenyl-d14	0.602	0.627	-4.2	118	0.00
33		3,3'-Dichlorobenzidine	0.336	0.361	-7.4	148	0.00
34	I	Perylene-d12	1.000	1.000	0.0	115	0.00

(#) = Out of Range SPCC's out = 1 CCC's out = 0