Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_M\Data\BM041015\
Data File : BM000950.D
Acq On : 10 Apr 2015 21:36
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Apr 11 00:10:10 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIM-BM040915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 10 23:03:27 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	167#	0.00
2		1,4-Dioxane	0.497	0.459	7.6	160#	0.00
3		n-Nitrosodimethylamine	0.486	0.487	-0.2	171#	0.00
4	S	2-Fluorophenol	1.058	1.059	-0.1	168#	0.00
5	S	Phenol-d6	1.267	1.262	0.4	171#	0.00
6	C	Phenol	1.302	1.300	0.2	172#	-0.02
7		bis(2-Chloroethyl)ether	1.142	1.132	0.9	167#	0.00
8	I	Naphthalene-d8	1.000	1.000	0.0	165#	0.00
9	S	Nitrobenzene-d5	0.241	0.238	1.2	171#	0.00
10		Nitrobenzene	0.283	0.288	-1.8	175#	0.00
11		2,4-Dimethylphenol	0.275	0.277	-0.7	172#	0.00
12	C	2,4-Dichlorophenol	0.243	0.236	2.9	171#	0.00
13		Naphthalene	1.009	0.984	2.5	165#	0.00
14	C	Hexachlorobutadiene	0.184	0.175	4.9	162#	0.00
15	C	2-Methylnaphthalene	0.689	0.688	0.1	169#	0.00
16		1-Methylnaphthalene	0.694	0.693	0.1	169#	-0.01
17	I	Acenaphthene-d10	1.000	1.000	0.0	209#	0.01
18	S	2,4,6-Tribromophenol	0.179	0.150	16.2	142	0.01
19	S	2-Fluorobiphenyl	1.401	1.316	6.1	168#	0.00
20		Acenaphthylene	2.032	1.795	11.7	139	0.01
21	C	Acenaphthene	1.380	1.308	5.2	184#	-0.02
22	P	2,4-Dinitrophenol	0.058	0.041#	29.3#	210#	-0.01
23		Fluorene	1.606	1.630	-1.5	202#	-0.02
24	I	Phenanthrene-d10	1.000	1.000	0.0	176#	0.01
25		Hexachlorobenzene	0.270	0.258	4.4	160#	0.01
26	C	Pentachlorophenol	0.086	0.075	12.8	171#	0.00
27		Phenanthrene	1.292	1.259	2.6	162#	0.01
28		Anthracene	1.184	1.178	0.5	181#	-0.01
29	C	Fluoranthene	1.350	1.293	4.2	163#	-0.01
30	I	Chrysene-d12	1.000	1.000	0.0	168#	0.00
31		Benzidine	0.300	0.253	15.7	172#	0.00
32		Pyrene	1.625	1.596	1.8	163#	0.00
33	S	Terphenyl-d14	0.609	0.639	-4.9	180#	0.00
34		Benzo(a)anthracene	1.443	1.415	1.9	169#	0.00
35		3,3'-Dichlorobenzidine	0.389	0.385	1.0	183#	0.00
36		Chrysene	1.475	1.444	2.1	163#	0.00
37		Indeno(1,2,3-cd)pyrene	1.506	1.493	0.9	168#	0.00
38	I	Perylene-d12	1.000	1.000	0.0	167#	0.00
39		Benzo(b)fluoranthene	1.565	1.520	2.9	171#	0.00
40		Benzo(k)fluoranthene	1.506	1.483	1.5	161#	0.00
41	C	Benzo(a)pyrene	1.378	1.344	2.5	167#	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Dibenzo(a,h)anthracene	1.278	1.273	0.4	169#	0.00
43	Benzo(g,h,i)perylene	1.410	1.382	2.0	165#	0.00

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