Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM041315\
Data File : BM000952.D
Acq On : 13 Apr 2015 16:23
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Apr 14 06:19:09 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 16:15:17 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	75	0.00
2		1,4-Dioxane	1.000	1.114	-11.4	80	0.00
3		n-Nitrosodimethylamine	1.000	0.991	0.9	75	0.00
4	S	2-Fluorophenol	1.000	1.010	-1.0	75	-0.02
5	S	Phenol-d6	1.000	0.966	3.4	74	0.00
6	C	Phenol	1.000	0.960	4.0	74	-0.02
7		bis(2-Chloroethyl)ether	1.000	0.990	1.0	74	0.00
8	I	Naphthalene-d8	5.000	5.000	0.0	72	0.00
9	S	Nitrobenzene-d5	1.000	0.988	1.2	75	0.00
10		Nitrobenzene	1.000	0.956	4.4	75	0.00
11		2,4-Dimethylphenol	1.000	0.961	3.9	72	0.00
12	C	2,4-Dichlorophenol	1.000	0.921	7.9	71	0.00
13		Naphthalene	1.000	1.029	-2.9	76	0.00
14	C	Hexachlorobutadiene	1.000	1.033	-3.3	77	0.00
15	C	2-Methylnaphthalene	1.000	1.033	-3.3	77	0.00
16		1-Methylnaphthalene	1.000	1.034	-3.4	77	-0.01
17	I	Acenaphthene-d10	5.000	5.000	0.0	74	0.00
18	S	2,4,6-Tribromophenol	1.000	0.998	0.2	61	0.00
19	S	2-Fluorobiphenyl	1.000	1.206	-20.6	76	0.00
20		Acenaphthylene	1.000	1.202	-20.2	70	0.00
21	C	Acenaphthene	1.000	1.033	-3.3	71	0.00
22	P	2,4-Dinitrophenol	1.000	0.763	23.7	55	0.00
23		Fluorene	1.000	0.987	1.3	70	0.00
24	I	Phenanthrene-d10	5.000	5.000	0.0	67	0.00
25		Hexachlorobenzene	1.000	1.118	-11.8	72	0.00
26	C	Pentachlorophenol	1.000	0.823	17.7	62	0.00
27		Phenanthrene	1.000	1.110	-11.0	70	0.00
28		Anthracene	1.000	0.975	2.5	68	0.00
29	C	Fluoranthene	1.000	0.993	0.7	65	0.00
30	I	Chrysene-d12	5.000	5.000	0.0	62	0.00
31		Benzidine	1.000	0.784	21.6	50	0.02
32		Pyrene	1.000	1.041	-4.1	63	0.00
33	S	Terphenyl-d14	1.000	1.026	-2.6	65	0.02
34		Benzo(a)anthracene	1.000	1.054	-5.4	67	0.00
35		3,3'-Dichlorobenzidine	1.000	0.839	16.1	55	0.02
36		Chrysene	1.000	1.005	-0.5	61	0.00
37		Indeno(1,2,3-cd)pyrene	1.000	0.996	0.4	62	0.00
38	I	Perylene-d12	5.000	5.000	0.0	60	0.00
39		Benzo(b)fluoranthene	1.000	1.108	-10.8	70	0.01
40		Benzo(k)fluoranthene	1.000	1.002	-0.2	58	0.01
41	C	Benzo(a)pyrene	1.000	1.011	-1.1	62	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM041315\
Data File : BM000952.D
Acq On : 13 Apr 2015 16:23
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Apr 14 06:19:09 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 16:15:17 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	Amount	Calc.	%Dev	Area%	Dev(min)
42	Dibenzo(a,h)anthracene	1.000	1.013	-1.3	61	0.00
43	Benzo(g,h,i)perylene	1.000	1.059	-5.9	64	0.00

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