Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM040915\
Data File : BM000935.D
Acq On : 09 Apr 2015 23:09
Operator : TP/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM040915

Quant Time: Apr 10 16:34:28 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	103	0.00
2		1,4-Dioxane	1.000	1.072	-7.2	106	0.00
3		n-Nitrosodimethylamine	1.000	1.025	-2.5	108	0.00
4	S	2-Fluorophenol	1.000	1.063	-6.3	109	0.00
5	S	Phenol-d6	1.000	1.060	-6.0	112	0.00
6	C	Phenol	1.000	1.050	-5.0	111	0.00
7		bis(2-Chloroethyl)ether	1.000	1.064	-6.4	110	0.00
8	I	Naphthalene-d8	5.000	5.000	0.0	99	0.00
9	S	Nitrobenzene-d5	1.000	1.075	-7.5	112	0.00
10		Nitrobenzene	1.000	1.027	-2.7	112	0.00
11		2,4-Dimethylphenol	1.000	1.069	-6.9	110	0.00
12	C	2,4-Dichlorophenol	1.000	1.059	-5.9	112	0.00
13		Naphthalene	1.000	1.048	-4.8	107	0.00
14	C	Hexachlorobutadiene	1.000	1.067	-6.7	109	0.00
15	C	2-Methylnaphthalene	1.000	1.072	-7.2	109	0.00
16		1-Methylnaphthalene	1.000	1.082	-8.2	110	-0.01
17	I	Acenaphthene-d10	5.000	5.000	0.0	125	0.01
18	S	2,4,6-Tribromophenol	1.000	0.908	9.2	93	0.01
19	S	2-Fluorobiphenyl	1.000	1.026	-2.6	110	0.00
20		Acenaphthylene	1.000	0.898	10.2	88	0.01
21	C	Acenaphthene	1.000	1.058	-5.8	123	-0.02
22	P	2,4-Dinitrophenol	1.000	1.015	-1.5	140	-0.01
23		Fluorene	1.000	1.127	-12.7	134	-0.02
24	I	Phenanthrene-d10	5.000	5.000	0.0	107	0.01
25		Hexachlorobenzene	1.000	1.026	-2.6	105	0.01
26	C	Pentachlorophenol	1.000	0.900	10.0	111	0.00
27		Phenanthrene	1.000	1.026	-2.6	104	0.01
28		Anthracene	1.000	1.055	-5.5	117	-0.01
29	C	Fluoranthene	1.000	1.043	-4.3	108	-0.01
30	I	Chrysene-d12	5.000	5.000	0.0	102	0.00
31		Benzidine	1.000	1.095	-9.5	121	0.00
32		Pyrene	1.000	1.069	-6.9	108	0.00
33	S	Terphenyl-d14	1.000	1.098	-9.8	114	0.00
34		Benzo(a)anthracene	1.000	1.071	-7.1	113	-0.01
35		3,3'-Dichlorobenzidine	1.000	1.004	-0.4	119	0.00
36		Chrysene	1.000	1.054	-5.4	106	0.00
37		Indeno(1,2,3-cd)pyrene	1.000	1.090	-9.0	112	0.00
38	I	Perylene-d12	5.000	5.000	0.0	102	0.00
39		Benzo(b)fluoranthene	1.000	1.120	-12.0	120	0.00
40		Benzo(k)fluoranthene	1.000	1.018	-1.8	101	0.00
41	C	Benzo(a)pyrene	1.000	1.065	-6.5	111	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM040915\
Data File : BM000935.D
Acq On : 09 Apr 2015 23:09
Operator : TP/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM040915

Quant Time: Apr 10 16:34:28 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	1.105	-10.5	114	0.00
43	Benzo(g,h,i)perylene	1.000	1.073	-7.3	110	0.00

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