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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	75	0.00
2		1,4-Dioxane	0.497	0.517	-4.0	80	0.00
3		n-Nitrosodimethylamine	0.486	0.482	0.8	75	0.00
4	S	2-Fluorophenol	1.058	1.069	-1.0	75	-0.02
5	S	Phenol-d6	1.267	1.223	3.5	74	0.00
6	C	Phenol	1.302	1.250	4.0	74	-0.02
7		bis(2-Chloroethyl)ether	1.142	1.131	1.0	74	0.00
8	I	Naphthalene-d8	1.000	1.000	0.0	72	0.00
9	S	Nitrobenzene-d5	0.241	0.239	0.8	75	0.00
10		Nitrobenzene	0.283	0.283	0.0	75	0.00
11		2,4-Dimethylphenol	0.275	0.265	3.6	72	0.00
12	C	2,4-Dichlorophenol	0.243	0.224	7.8	71	0.00
13		Naphthalene	1.009	1.039	-3.0	76	0.00
14	C	Hexachlorobutadiene	0.184	0.191	-3.8	77	0.00
15	C	2-Methylnaphthalene	0.689	0.711	-3.2	77	0.00
16		1-Methylnaphthalene	0.694	0.718	-3.5	77	-0.01
17	I	Acenaphthene-d10	1.000	1.000	0.0	74	0.00
18	S	2,4,6-Tribromophenol	0.179	0.183	-2.2	61	0.00
19	S	2-Fluorobiphenyl	1.401	1.689	-20.6	76	0.00
20		Acenaphthylene	2.032	2.543	-25.1#	70	0.00
21	C	Acenaphthene	1.380	1.425	-3.3	71	0.00
22	P	2,4-Dinitrophenol	0.058	0.031#	46.6#	55	0.00
23		Fluorene	1.606	1.585	1.3	70	0.00
24	I	Phenanthrene-d10	1.000	1.000	0.0	67	0.00
25		Hexachlorobenzene	0.270	0.302	-11.9	72	0.00
26	C	Pentachlorophenol	0.086	0.071	17.4	62	0.00
27		Phenanthrene	1.292	1.434	-11.0	70	0.00
28		Anthracene	1.184	1.155	2.4	68	0.00
29	C	Fluoranthene	1.350	1.341	0.7	65	0.00
30	I	Chrysene-d12	1.000	1.000	0.0	62	0.00
31		Benzidine	0.300	0.202	32.7#	50	0.02
32		Pyrene	1.625	1.692	-4.1	63	0.00
33	S	Terphenyl-d14	0.609	0.625	-2.6	65	0.02
34		Benzo(a)anthracene	1.443	1.521	-5.4	67	0.00
35		3,3'-Dichlorobenzidine	0.389	0.315	19.0	55	0.02
36		Chrysene	1.475	1.483	-0.5	61	0.00
37		Indeno(1,2,3-cd)pyrene	1.506	1.500	0.4	62	0.00
38	I	Perylene-d12	1.000	1.000	0.0	60	0.00
39		Benzo(b)fluoranthene	1.565	1.734	-10.8	70	0.01
40		Benzo(k)fluoranthene	1.506	1.508	-0.1	58	0.01
41	C	Benzo(a)pyrene	1.378	1.393	-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	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Dibenzo(a,h)anthracene	1.278	1.294	-1.3	61	0.00
43	Benzo(g,h,i)perylene	1.410	1.493	-5.9	64	0.00

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