Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM111615\
Data File : BM003136.D
Acq On : 16 Nov 2015 10:47
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Nov 17 13:42:35 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM110415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 17 12:50:51 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	88	-0.02
2		1,4-Dioxane	1.000	0.965	3.5	83	-0.02
3		n-Nitrosodimethylamine	1.000	0.938	6.2	85	-0.02
4	S	2-Fluorophenol	1.000	0.903	9.7	83	-0.02
5	S	Phenol-d6	1.000	0.897	10.3	83	-0.02
6		bis(2-Chloroethyl)ether	1.000	0.951	4.9	86	-0.03
7	I	Naphthalene-d8	5.000	5.000	0.0	87	-0.01
8	S	Nitrobenzene-d5	1.000	0.957	4.3	88	-0.02
9		Nitrobenzene	1.000	0.953	4.7	88	-0.02
10		Naphthalene	1.000	0.990	1.0	88	-0.01
11		Hexachlorobutadiene	1.000	1.000	0.0	88	-0.01
12		2-Methylnaphthalene	1.000	0.979	2.1	87	-0.02
13	I	Acenaphthene-d10	5.000	5.000	0.0	84	0.00
14	S	2,4,6-Tribromophenol	1.000	0.975	2.5	83	0.00
15	S	2-Fluorobiphenyl	1.000	1.034	-3.4	88	-0.02
16		Acenaphthylene	1.000	0.940	6.0	82	0.00
17		Acenaphthene	1.000	1.053	-5.3	96	-0.03
18		Fluorene	1.000	1.213	-21.3	100	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	88	0.00
20		4-Bromophenyl-phenylether	1.000	1.766	-76.6#	106	0.00
21		Hexachlorobenzene	1.000	1.310	-31.0#	83	0.00
22		Pentachlorophenol	1.000	1.275	-27.5#	93	0.00
23		Phenanthrene	1.000	1.799	-79.9#	103	0.00
24		Anthracene	1.000	0.971	2.9	85	0.00
25		Fluoranthene	1.000	1.300	-30.0#	96	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	101	0.00
27		Benzidine	1.000	1.096	-9.6	97	0.00
28		Pyrene	1.000	0.941	5.9	97	0.00
29	S	Terphenyl-d14	1.000	0.912	8.8	89	0.00
30		Benzo(a)anthracene	1.000	1.007	-0.7	102	0.00
31		3,3'-Dichlorobenzidine	1.000	1.025	-2.5	91	0.00
32		Chrysene	1.000	0.934	6.6	82	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	0.757	24.3	74	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	1.031	-3.1	92	-0.02
35	I	Perylene-d12	5.000	5.000	0.0	92	0.00
36		Benzo(b)fluoranthene	1.000	0.889	11.1	84	0.00
37		Benzo(k)fluoranthene	1.000	0.973	2.7	91	0.00
38	C	Benzo(a)pyrene	1.000	0.906	9.4	92	0.00
39		Dibenzo(a,h)anthracene	1.000	0.956	4.4	93	0.00
40		Benzo(g,h,i)perylene	1.000	0.922	7.8	91	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM111615\
Data File : BM003136.D
Acq On : 16 Nov 2015 10:47
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Nov 17 13:42:35 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM110415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 17 12:50:51 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)

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