Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122415\
Data File : BM003772.D
Acq On : 24 Dec 2015 10:49
Operator : UM/SJ
Sample : SSTDCCC001
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Dec 24 16:10:58 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM122415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 24 10:08:49 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	0.00
2		1,4-Dioxane	1.000	0.915	8.5	89	0.00
3		n-Nitrosodimethylamine	1.000	0.828	17.2	84	0.00
4	S	2-Fluorophenol	1.000	0.829	17.1	86	0.02
5	S	Phenol-d6	1.000	0.807	19.3	86	0.00
6		bis(2-Chloroethyl)ether	1.000	0.832	16.8	85	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	89	0.00
8	S	Nitrobenzene-d5	1.000	0.806	19.4	85	0.00
9		Nitrobenzene	1.000	0.802	19.8	84	0.00
10		Naphthalene	1.000	0.880	12.0	89	0.00
11		Hexachlorobutadiene	1.000	0.884	11.6	89	0.00
12		2-Methylnaphthalene	1.000	0.867	13.3	88	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	91	0.00
14	S	2,4,6-Tribromophenol	1.000	0.836	16.4	91	0.00
15	S	2-Fluorobiphenyl	1.000	0.852	14.8	89	0.00
16		Acenaphthylene	1.000	0.799	20.1	88	0.00
17		Acenaphthene	1.000	0.919	8.1	91	0.00
18		Fluorene	1.000	0.860	14.0	90	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	97	0.03
20		4-Bromophenyl-phenylether	1.000	1.132	-13.2	88	0.00
21		Hexachlorobenzene	1.000	1.194	-19.4	93	0.00
22		Pentachlorophenol	1.000	0.874	12.6	107	0.00
23		Phenanthrene	1.000	1.142	-14.2	96	0.00
24		Anthracene	1.000	0.905	9.5	95	0.00
25		Fluoranthene	1.000	0.959	4.1	98	0.02
26	I	Chrysene-d12	5.000	5.000	0.0	120	0.00
27		Benzidine	1.000	0.895	10.5	105	0.00
28		Pyrene	1.000	0.769	23.1	101	0.00
29	S	Terphenyl-d14	1.000	0.794	20.6	103	0.00
30		Benzo(a)anthracene	1.000	0.806	19.4	116	0.00
31		3,3'-Dichlorobenzidine	1.000	0.894	10.6	108	0.00
32		Chrysene	1.000	0.846	15.4	107	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	0.776	22.4	100	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	0.972	2.8	125	0.01
35	I	Perylene-d12	5.000	5.000	0.0	125	0.01
36		Benzo(b)fluoranthene	1.000	0.836	16.4	116	0.00
37		Benzo(k)fluoranthene	1.000	0.846	15.4	127	0.00
38	C	Benzo(a)pyrene	1.000	0.836	16.4	123	0.00
39		Dibenzo(a,h)anthracene	1.000	0.878	12.2	125	0.00
40		Benzo(g,h,i)perylene	1.000	0.877	12.3	124	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122415\
Data File : BM003772.D
Acq On : 24 Dec 2015 10:49
Operator : UM/SJ
Sample : SSTDCCC001
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC001

Quant Time: Dec 24 16:10:58 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM122415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 24 10:08:49 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