Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM100815\
Data File : BM002878.D
Acq On : 09 Oct 2015 00:00
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Oct 09 07:12:29 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM100815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Oct 09 06:47:52 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	65	0.00
2		1,4-Dioxane	1.000	1.008	-0.8	64	0.00
3		n-Nitrosodimethylamine	1.000	0.976	2.4	62	0.00
4	S	2-Fluorophenol	1.000	0.964	3.6	64	0.02
5	S	Phenol-d6	1.000	0.910	9.0	62	0.02
6		bis(2-Chloroethyl)ether	1.000	0.979	2.1	66	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	67	0.00
8	S	Nitrobenzene-d5	1.000	0.925	7.5	63	0.01
9		Nitrobenzene	1.000	0.910	9.0	63	0.00
10		Naphthalene	1.000	0.982	1.8	68	0.00
11		Hexachlorobutadiene	1.000	0.984	1.6	68	0.00
12		2-Methylnaphthalene	1.000	1.004	-0.4	69	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	71	0.00
14	S	2,4,6-Tribromophenol	1.000	0.987	1.3	75	0.00
15	S	2-Fluorobiphenyl	1.000	0.990	1.0	70	0.00
16		Acenaphthylene	1.000	0.938	6.2	67	0.00
17		Acenaphthene	1.000	0.947	5.3	70	0.00
18		Fluorene	1.000	0.986	1.4	71	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	79	0.00
20		4-Bromophenyl-phenylether	1.000	0.958	4.2	74	0.00
21		Hexachlorobenzene	1.000	0.905	9.5	73	0.00
22		Pentachlorophenol	1.000	0.966	3.4	69	0.00
23		Phenanthrene	1.000	0.936	6.4	76	0.00
24		Anthracene	1.000	0.909	9.1	75	0.00
25		Fluoranthene	1.000	0.994	0.6	84	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	90	0.00
27		Benzidine	1.000	0.895	10.5	88	0.00
28		Pyrene	1.000	0.944	5.6	83	0.00
29	S	Terphenyl-d14	1.000	0.965	3.5	89	-0.02
30		Benzo(a)anthracene	1.000	0.946	5.4	91	0.00
31		3,3'-Dichlorobenzidine	1.000	0.948	5.2	89	0.00
32		Chrysene	1.000	1.034	-3.4	94	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	0.810	19.0	62	-0.02
34		Indeno(1,2,3-cd)pyrene	1.000	1.097	-9.7	102	0.00
35	I	Perylene-d12	5.000	5.000	0.0	98	0.00
36		Benzo(b)fluoranthene	1.000	1.018	-1.8	98	0.00
37		Benzo(k)fluoranthene	1.000	0.928	7.2	97	0.00
38	C	Benzo(a)pyrene	1.000	0.964	3.6	97	-0.01
39		Dibenzo(a,h)anthracene	1.000	1.007	-0.7	102	0.00
40		Benzo(g,h,i)perylene	1.000	1.012	-1.2	102	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM100815\
Data File : BM002878.D
Acq On : 09 Oct 2015 00:00
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Oct 09 07:12:29 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM100815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Oct 09 06:47:52 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