Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM010916\
Data File : BM003933.D
Acq On : 10 Jan 2016 11:19
Operator : SJ/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Jan 11 00:59:50 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM010216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 04 13:33:21 2016
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	71	0.00
2		1,4-Dioxane	1.000	1.021	-2.1	72	0.00
3		n-Nitrosodimethylamine	1.000	0.849	15.1	60	0.00
4	S	2-Fluorophenol	1.000	0.895	10.5	67	0.00
5	S	Phenol-d6	1.000	0.901	9.9	65	0.00
6		bis(2-Chloroethyl)ether	1.000	0.944	5.6	66	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	69	0.00
8	S	Nitrobenzene-d5	1.000	0.835	16.5	58	0.00
9		Nitrobenzene	1.000	0.885	11.5	62	0.00
10		Naphthalene	1.000	0.978	2.2	68	0.00
11		Hexachlorobutadiene	1.000	0.966	3.4	68	0.00
12		2-Methylnaphthalene	1.000	0.904	9.6	62	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	66	0.00
14	S	2,4,6-Tribromophenol	1.000	0.816	18.4	55	0.00
15	S	2-Fluorobiphenyl	1.000	0.933	6.7	62	0.00
16		Acenaphthylene	1.000	0.917	8.3	60	0.00
17		Acenaphthene	1.000	0.929	7.1	65	0.00
18		Fluorene	1.000	0.966	3.4	63	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	64	0.00
20		4-Bromophenyl-phenylether	1.000	0.975	2.5	67	0.00
21		Hexachlorobenzene	1.000	0.930	7.0	66	0.00
22		Pentachlorophenol	1.000	0.745	25.5#	47	0.00
23		Phenanthrene	1.000	1.046	-4.6	66	0.00
24		Anthracene	1.000	0.923	7.7	64	0.00
25		Fluoranthene	1.000	0.951	4.9	63	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	71	0.00
27		Benzidine	1.000	1.173	-17.3	85	0.00
28		Pyrene	1.000	0.912	8.8	64	0.00
29	S	Terphenyl-d14	1.000	0.935	6.5	67	0.00
30		Benzo(a)anthracene	1.000	0.906	9.4	66	0.02
31		3,3'-Dichlorobenzidine	1.000	0.867	13.3	63	0.00
32		Chrysene	1.000	1.106	-10.6	77	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	0.904	9.6	71	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	0.924	7.6	66	0.01
35	I	Perylene-d12	5.000	5.000	0.0	76	0.00
36		Benzo(b)fluoranthene	1.000	0.995	0.5	81	0.00
37		Benzo(k)fluoranthene	1.000	0.857	14.3	66	0.01
38	C	Benzo(a)pyrene	1.000	0.894	10.6	72	0.00
39		Dibenzo(a,h)anthracene	1.000	0.860	14.0	65	0.00
40		Benzo(g,h,i)perylene	1.000	0.885	11.5	68	0.01

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

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