Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM010916\
Data File : BM003931.D
Acq On : 10 Jan 2016 09:28
Operator : SJ/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Jan 11 00:49:57 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	80	0.00
2		1,4-Dioxane	1.000	1.045	-4.5	83	0.00
3		n-Nitrosodimethylamine	1.000	0.848	15.2	68	0.00
4	S	2-Fluorophenol	1.000	0.931	6.9	79	0.00
5	S	Phenol-d6	1.000	0.907	9.3	74	0.00
6		bis(2-Chloroethyl)ether	1.000	0.947	5.3	74	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	73	0.00
8	S	Nitrobenzene-d5	1.000	0.858	14.2	64	0.00
9		Nitrobenzene	1.000	0.907	9.3	68	-0.01
10		Naphthalene	1.000	0.961	3.9	71	-0.02
11		Hexachlorobutadiene	1.000	0.968	3.2	73	0.00
12		2-Methylnaphthalene	1.000	0.859	14.1	63	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	63	0.00
14	S	2,4,6-Tribromophenol	1.000	0.818	18.2	52	0.00
15	S	2-Fluorobiphenyl	1.000	0.972	2.8	61	0.00
16		Acenaphthylene	1.000	0.925	7.5	57	0.00
17		Acenaphthene	1.000	0.979	2.1	65	0.00
18		Fluorene	1.000	0.950	5.0	59	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	61	0.00
20		4-Bromophenyl-phenylether	1.000	0.903	9.7	58	0.00
21		Hexachlorobenzene	1.000	0.894	10.6	60	0.00
22		Pentachlorophenol	1.000	0.876	12.4	55	0.00
23		Phenanthrene	1.000	0.980	2.0	58	0.00
24		Anthracene	1.000	0.940	6.0	61	0.00
25		Fluoranthene	1.000	0.917	8.3	58	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	74	0.00
27		Benzidine	1.000	1.204	-20.4	91	0.00
28		Pyrene	1.000	0.790	21.0	58	0.00
29	S	Terphenyl-d14	1.000	0.788	21.2	58	0.00
30		Benzo(a)anthracene	1.000	0.895	10.5	68	0.00
31		3,3'-Dichlorobenzidine	1.000	0.886	11.4	67	0.00
32		Chrysene	1.000	0.974	2.6	71	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	0.822	17.8	63	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	1.219	-21.9	90	0.01
35	I	Perylene-d12	5.000	5.000	0.0	89	0.00
36		Benzo(b)fluoranthene	1.000	0.860	14.0	82	0.00
37		Benzo(k)fluoranthene	1.000	0.901	9.9	82	0.00
38	C	Benzo(a)pyrene	1.000	0.911	8.9	86	0.00
39		Dibenzo(a,h)anthracene	1.000	1.010	-1.0	91	0.00
40		Benzo(g,h,i)perylene	1.000	1.025	-2.5	93	0.01

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

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