Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM120415\
Data File : BM003373.D
Acq On : 03 Dec 2015 16:14
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Dec 04 00:22:28 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM120215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 02 16:19:00 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	70	0.00
2		1,4-Dioxane	1.000	0.935	6.5	74	0.00
3		n-Nitrosodimethylamine	1.000	0.996	0.4	73	-0.02
4	S	2-Fluorophenol	1.000	0.916	8.4	72	-0.02
5	S	Phenol-d6	1.000	0.927	7.3	69	0.00
6		bis(2-Chloroethyl)ether	1.000	0.951	4.9	70	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	68	0.00
8	S	Nitrobenzene-d5	1.000	0.954	4.6	70	0.00
9		Nitrobenzene	1.000	0.954	4.6	68	0.00
10		Naphthalene	1.000	0.952	4.8	70	0.00
11		Hexachlorobutadiene	1.000	0.997	0.3	73	0.00
12		2-Methylnaphthalene	1.000	0.946	5.4	67	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	67	0.00
14	S	2,4,6-Tribromophenol	1.000	0.952	4.8	63	0.00
15	S	2-Fluorobiphenyl	1.000	0.914	8.6	68	0.00
16		Acenaphthylene	1.000	0.945	5.5	65	0.00
17		Acenaphthene	1.000	0.888	11.2	67	0.00
18		Fluorene	1.000	0.917	8.3	65	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	66	0.00
20		4-Bromophenyl-phenylether	1.000	0.906	9.4	69	0.00
21		Hexachlorobenzene	1.000	0.894	10.6	65	0.00
22		Pentachlorophenol	1.000	0.935	6.5	63	0.00
23		Phenanthrene	1.000	0.906	9.4	68	0.00
24		Anthracene	1.000	0.890	11.0	60	0.00
25		Fluoranthene	1.000	0.974	2.6	64	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	68	0.00
27		Benzidine	1.000	0.854	14.6	60	0.02
28		Pyrene	1.000	0.839	16.1	62	0.00
29	S	Terphenyl-d14	1.000	0.835	16.5	60	0.00
30		Benzo(a)anthracene	1.000	0.920	8.0	65	0.00
31		3,3'-Dichlorobenzidine	1.000	0.774	22.6	57	0.00
32		Chrysene	1.000	0.770	23.0	55	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	0.683	31.7#	40	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	0.743	25.7#	65	0.01
35	I	Perylene-d12	5.000	5.000	0.0	60	0.01
36		Benzo(b)fluoranthene	1.000	0.974	2.6	61	0.00
37		Benzo(k)fluoranthene	1.000	0.922	7.8	57	0.01
38	C	Benzo(a)pyrene	1.000	0.905	9.5	58	0.01
39		Dibenzo(a,h)anthracene	1.000	0.900	10.0	67	0.01
40		Benzo(g,h,i)perylene	1.000	0.879	12.1	67	0.01

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

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