Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM010816\
Data File : BM003924.D
Acq On : 08 Jan 2016 16:10
Operator : SJ/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Jan 09 00:28:19 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	89	0.00
2		1,4-Dioxane	1.000	0.991	0.9	88	0.00
3		n-Nitrosodimethylamine	1.000	0.858	14.2	77	0.00
4	S	2-Fluorophenol	1.000	0.938	6.2	89	0.00
5	S	Phenol-d6	1.000	0.931	6.9	85	0.00
6		bis(2-Chloroethyl)ether	1.000	0.965	3.5	85	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	85	0.00
8	S	Nitrobenzene-d5	1.000	0.869	13.1	75	0.00
9		Nitrobenzene	1.000	0.920	8.0	80	0.00
10		Naphthalene	1.000	0.968	3.2	83	-0.02
11		Hexachlorobutadiene	1.000	0.972	2.8	85	0.00
12		2-Methylnaphthalene	1.000	0.893	10.7	76	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	78	0.00
14	S	2,4,6-Tribromophenol	1.000	0.875	12.5	70	0.00
15	S	2-Fluorobiphenyl	1.000	0.963	3.7	75	0.00
16		Acenaphthylene	1.000	0.966	3.4	74	0.00
17		Acenaphthene	1.000	0.992	0.8	82	0.00
18		Fluorene	1.000	1.001	-0.1	77	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	80	0.00
20		4-Bromophenyl-phenylether	1.000	0.883	11.7	75	0.00
21		Hexachlorobenzene	1.000	0.864	13.6	76	0.00
22		Pentachlorophenol	1.000	0.789	21.1	63	0.00
23		Phenanthrene	1.000	0.947	5.3	74	0.00
24		Anthracene	1.000	0.881	11.9	75	0.00
25		Fluoranthene	1.000	0.872	12.8	72	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	82	0.00
27		Benzidine	1.000	0.827	17.3	64	0.00
28		Pyrene	1.000	0.877	12.3	72	0.00
29	S	Terphenyl-d14	1.000	0.853	14.7	71	-0.02
30		Benzo(a)anthracene	1.000	0.973	2.7	83	0.00
31		3,3'-Dichlorobenzidine	1.000	0.778	22.2	63	0.00
32		Chrysene	1.000	0.760	24.0	62	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	0.741	25.9#	58	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	0.998	0.2	82	0.00
35	I	Perylene-d12	5.000	5.000	0.0	82	-0.01
36		Benzo(b)fluoranthene	1.000	0.932	6.8	82	0.00
37		Benzo(k)fluoranthene	1.000	0.881	11.9	74	0.00
38	C	Benzo(a)pyrene	1.000	0.937	6.3	81	0.00
39		Dibenzo(a,h)anthracene	1.000	1.000	0.0	82	-0.01
40		Benzo(g,h,i)perylene	1.000	1.024	-2.4	85	0.00

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

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