Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM012216\
Data File : BM004110.D
Acq On : 22 Jan 2016 14:39
Operator : SJ/UM
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM012216

Quant Time: Jan 22 15:17:53 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM012216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 22 14:46:41 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	107	0.00
2		1,4-Dioxane	1.000	0.908	9.2	102	0.00
3		n-Nitrosodimethylamine	1.000	0.898	10.2	105	-0.02
4	S	2-Fluorophenol	1.000	0.877	12.3	104	0.00
5	S	Phenol-d6	1.000	0.870	13.0	104	0.00
6		bis(2-Chloroethyl)ether	1.000	0.877	12.3	104	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	106	0.00
8	S	Nitrobenzene-d5	1.000	0.903	9.7	106	0.00
9		Nitrobenzene	1.000	0.898	10.2	106	0.00
10		Naphthalene	1.000	0.883	11.7	102	0.00
11		Hexachlorobutadiene	1.000	0.895	10.5	104	0.00
12		2-Methylnaphthalene	1.000	0.880	12.0	102	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	107	0.00
14	S	2,4,6-Tribromophenol	1.000	0.874	12.6	101	0.00
15	S	2-Fluorobiphenyl	1.000	0.877	12.3	102	0.00
16		Acenaphthylene	1.000	0.868	13.2	102	0.00
17		Acenaphthene	1.000	0.841	15.9	101	0.00
18		Fluorene	1.000	0.858	14.2	102	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	106	0.00
20		4-Bromophenyl-phenylether	1.000	0.920	8.0	100	0.00
21		Hexachlorobenzene	1.000	0.925	7.5	103	0.00
22		Pentachlorophenol	1.000	0.856	14.4	109	0.00
23		Phenanthrene	1.000	0.882	11.8	102	0.00
24		Anthracene	1.000	0.899	10.1	102	0.00
25		Fluoranthene	1.000	0.894	10.6	102	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	100	0.00
27		Benzidine	1.000	1.107	-10.7	115	0.00
28		Pyrene	1.000	0.933	6.7	102	0.00
29	S	Terphenyl-d14	1.000	0.940	6.0	104	0.00
30		Benzo(a)anthracene	1.000	0.860	14.0	98	0.00
31		3,3'-Dichlorobenzidine	1.000	1.035	-3.5	113	0.00
32		Chrysene	1.000	1.022	-2.2	114	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	1.031	-3.1	119	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	0.922	7.8	105	0.00
35	I	Perylene-d12	5.000	5.000	0.0	109	0.00
36		Benzo(b)fluoranthene	1.000	0.860	14.0	101	0.00
37		Benzo(k)fluoranthene	1.000	0.887	11.3	109	0.00
38	C	Benzo(a)pyrene	1.000	0.875	12.5	107	0.00
39		Dibenzo(a,h)anthracene	1.000	0.868	13.2	105	0.00
40		Benzo(g,h,i)perylene	1.000	0.851	14.9	104	0.00

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

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