Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM112715\
Data File : BM003309.D
Acq On : 27 Nov 2015 17:21
Operator : UM/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM112715

Quant Time: Nov 27 18:23:20 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM112715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 27 16:36:40 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	119	0.00
2		1,4-Dioxane	1.000	0.952	4.8	110	-0.02
3		n-Nitrosodimethylamine	1.000	0.984	1.6	117	0.00
4	S	2-Fluorophenol	1.000	0.978	2.2	120	0.00
5	S	Phenol-d6	1.000	1.009	-0.9	124	0.00
6		bis(2-Chloroethyl)ether	1.000	0.990	1.0	121	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	124	0.00
8	S	Nitrobenzene-d5	1.000	0.983	1.7	128	0.00
9		Nitrobenzene	1.000	0.984	1.6	130	0.00
10		Naphthalene	1.000	0.995	0.5	125	0.01
11		Hexachlorobutadiene	1.000	1.011	-1.1	129	0.00
12		2-Methylnaphthalene	1.000	1.037	-3.7	130	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	133	0.00
14	S	2,4,6-Tribromophenol	1.000	0.900	10.0	122	0.00
15	S	2-Fluorobiphenyl	1.000	0.926	7.4	133	0.00
16		Acenaphthylene	1.000	0.981	1.9	136	0.00
17		Acenaphthene	1.000	0.849	15.1	131	0.00
18		Fluorene	1.000	0.945	5.5	131	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	119	0.00
20		4-Bromophenyl-phenylether	1.000	0.891	10.9	127	-0.03
21		Hexachlorobenzene	1.000	0.998	0.2	130	0.00
22		Pentachlorophenol	1.000	0.808	19.2	110	0.00
23		Phenanthrene	1.000	0.880	12.0	119	0.00
24		Anthracene	1.000	1.002	-0.2	120	0.00
25		Fluoranthene	1.000	0.799	20.1	107	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	84	0.00
27		Benzidine	1.000	1.203	-20.3	123	0.00
28		Pyrene	1.000	1.075	-7.5	107	0.00
29	S	Terphenyl-d14	1.000	1.114	-11.4	102	0.00
30		Benzo(a)anthracene	1.000	1.035	-3.5	97	0.00
31		3,3'-Dichlorobenzidine	1.000	0.945	5.5	100	0.02
32		Chrysene	1.000	1.019	-1.9	95	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	1.147	-14.7	120	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	1.046	-4.6	101	0.01
35	I	Perylene-d12	5.000	5.000	0.0	92	0.01
36		Benzo(b)fluoranthene	1.000	0.972	2.8	88	0.01
37		Benzo(k)fluoranthene	1.000	1.044	-4.4	101	0.00
38	C	Benzo(a)pyrene	1.000	0.995	0.5	96	0.01
39		Dibenzo(a,h)anthracene	1.000	1.126	-12.6	105	0.01
40		Benzo(g,h,i)perylene	1.000	1.020	-2.0	97	0.01

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

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