Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM100815\
Data File : BM002874.D
Acq On : 08 Oct 2015 21:34
Operator : UM/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM100815

Quant Time: Oct 09 07:00:11 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM100815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Oct 09 06:47:52 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	105	0.00
2		1,4-Dioxane	1.000	0.950	5.0	97	0.00
3		n-Nitrosodimethylamine	1.000	0.939	6.1	96	0.00
4	S	2-Fluorophenol	1.000	0.910	9.0	97	0.02
5	S	Phenol-d6	1.000	0.888	11.2	96	0.00
6		bis(2-Chloroethyl)ether	1.000	0.896	10.4	97	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	105	0.00
8	S	Nitrobenzene-d5	1.000	0.927	7.3	99	0.01
9		Nitrobenzene	1.000	0.922	7.8	99	0.00
10		Naphthalene	1.000	0.911	8.9	99	0.00
11		Hexachlorobutadiene	1.000	0.915	8.5	99	0.00
12		2-Methylnaphthalene	1.000	0.914	8.6	98	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	105	0.00
14	S	2,4,6-Tribromophenol	1.000	0.862	13.8	97	0.00
15	S	2-Fluorobiphenyl	1.000	0.943	5.7	99	0.00
16		Acenaphthylene	1.000	0.918	8.2	97	0.00
17		Acenaphthene	1.000	0.885	11.5	97	0.00
18		Fluorene	1.000	0.906	9.4	97	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	104	0.00
20		4-Bromophenyl-phenylether	1.000	0.943	5.7	96	0.00
21		Hexachlorobenzene	1.000	0.903	9.7	96	0.00
22		Pentachlorophenol	1.000	0.900	10.0	84	0.00
23		Phenanthrene	1.000	0.911	8.9	97	0.00
24		Anthracene	1.000	0.887	11.3	96	0.00
25		Fluoranthene	1.000	0.865	13.5	96	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	99	0.00
27		Benzidine	1.000	0.916	8.4	99	0.00
28		Pyrene	1.000	0.987	1.3	96	0.00
29	S	Terphenyl-d14	1.000	0.964	3.6	97	0.00
30		Benzo(a)anthracene	1.000	0.875	12.5	93	0.00
31		3,3'-Dichlorobenzidine	1.000	0.897	10.3	92	0.00
32		Chrysene	1.000	0.967	3.3	96	-0.02
33		Bis(2-ethylhexyl)phthalate	1.000	0.916	8.4	77	-0.02
34		Indeno(1,2,3-cd)pyrene	1.000	0.895	10.5	91	0.00
35	I	Perylene-d12	5.000	5.000	0.0	99	0.00
36		Benzo(b)fluoranthene	1.000	0.907	9.3	88	0.00
37		Benzo(k)fluoranthene	1.000	0.942	5.8	100	0.00
38	C	Benzo(a)pyrene	1.000	0.905	9.5	93	0.00
39		Dibenzo(a,h)anthracene	1.000	0.888	11.2	91	0.00
40		Benzo(g,h,i)perylene	1.000	0.888	11.2	91	0.00

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

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