Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM110415\
Data File : BM003057.D
Acq On : 04 Nov 2015 23:26
Operator : UM/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : ICVBM110415

Quant Time: Nov 05 10:37:08 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-SIM-BM110415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Nov 05 10:29:18 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	183	0.00
2		1,4-Dioxane	1.000	0.903	9.7	163	0.00
3		n-Nitrosodimethylamine	1.000	0.918	8.2	175	0.00
4	S	2-Fluorophenol	1.000	0.878	12.2	168	0.00
5	S	Phenol-d6	1.000	0.942	5.8	183	0.00
6		bis(2-Chloroethyl)ether	1.000	0.956	4.4	180	-0.02
7	I	Naphthalene-d8	5.000	5.000	0.0	198	0.00
8	S	Nitrobenzene-d5	1.000	0.935	6.5	196	0.00
9		Nitrobenzene	1.000	0.946	5.4	198	0.00
10		Naphthalene	1.000	0.928	7.2	187	0.00
11		Hexachlorobutadiene	1.000	0.907	9.3	181	0.00
12		2-Methylnaphthalene	1.000	0.966	3.4	194	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	220	0.00
14	S	2,4,6-Tribromophenol	1.000	0.987	1.3	222	0.00
15	S	2-Fluorobiphenyl	1.000	0.884	11.6	197	0.00
16		Acenaphthylene	1.000	0.955	4.5	220	0.00
17		Acenaphthene	1.000	0.928	7.2	222	0.00
18		Fluorene	1.000	0.939	6.1	203	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	289	0.00
20		4-Bromophenyl-phenylether	1.000	0.867	13.3	174	0.00
21		Hexachlorobenzene	1.000	0.909	9.1	193	0.00
22		Pentachlorophenol	1.000	0.889	11.1	198	0.00
23		Phenanthrene	1.000	0.866	13.4	175	0.00
24		Anthracene	1.000	0.951	4.9	274	0.00
25		Fluoranthene	1.000	0.952	4.8	237	0.01
26	I	Chrysene-d12	5.000	5.000	0.0	257	0.00
27		Benzidine	1.000	0.991	0.9	226	0.00
28		Pyrene	1.000	0.939	6.1	247	0.00
29	S	Terphenyl-d14	1.000	0.922	7.8	230	0.00
30		Benzo(a)anthracene	1.000	0.959	4.1	248	0.00
31		3,3'-Dichlorobenzidine	1.000	0.936	6.4	208	0.00
32		Chrysene	1.000	0.825	17.5	189	0.01
33		Bis(2-ethylhexyl)phthalate	1.000	0.902	9.8	246	0.01
34		Indeno(1,2,3-cd)pyrene	1.000	0.802	19.8	179	0.00
35	I	Perylene-d12	5.000	5.000	0.0	207	-0.01
36		Benzo(b)fluoranthene	1.000	0.943	5.7	199	0.00
37		Benzo(k)fluoranthene	1.000	0.896	10.4	188	0.00
38	C	Benzo(a)pyrene	1.000	0.871	12.9	199	0.00
39		Dibenzo(a,h)anthracene	1.000	0.845	15.5	184	0.00
40		Benzo(g,h,i)perylene	1.000	0.807	19.3	179	0.00

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

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