Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM013016\
Data File : BM004147.D
Acq On : 29 Jan 2016 15:37
Operator : SJ/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : DFTPP

Quant Time: Jan 29 22:50:54 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	88	0.00
2		1,4-Dioxane	1.000	0.977	2.3	90	0.00
3		n-Nitrosodimethylamine	1.000	0.898	10.2	86	0.00
4	S	2-Fluorophenol	1.000	0.864	13.6	84	0.00
5	S	Phenol-d6	1.000	0.875	12.5	86	0.00
6		bis(2-Chloroethyl)ether	1.000	0.915	8.5	89	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	92	0.00
8	S	Nitrobenzene-d5	1.000	0.890	11.0	91	0.00
9		Nitrobenzene	1.000	0.893	10.7	91	0.00
10		Naphthalene	1.000	0.923	7.7	92	0.00
11		Hexachlorobutadiene	1.000	0.908	9.2	91	0.00
12		2-Methylnaphthalene	1.000	0.932	6.8	94	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	97	0.00
14	S	2,4,6-Tribromophenol	1.000	0.888	11.2	95	0.00
15	S	2-Fluorobiphenyl	1.000	0.901	9.9	95	0.00
16		Acenaphthylene	1.000	0.887	11.3	96	0.00
17		Acenaphthene	1.000	0.878	12.2	96	0.00
18		Fluorene	1.000	0.925	7.5	101	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	104	0.00
20		4-Bromophenyl-phenylether	1.000	0.894	10.6	95	0.00
21		Hexachlorobenzene	1.000	0.881	11.9	97	0.00
22		Pentachlorophenol	1.000	0.824	17.6	101	0.00
23		Phenanthrene	1.000	0.893	10.7	101	0.00
24		Anthracene	1.000	0.870	13.0	97	0.00
25		Fluoranthene	1.000	0.881	11.9	99	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	85	0.00
27		Benzidine	1.000	1.056	-5.6	92	0.02
28		Pyrene	1.000	1.047	-4.7	97	0.00
29	S	Terphenyl-d14	1.000	1.037	-3.7	98	0.00
30		Benzo(a)anthracene	1.000	0.877	12.3	84	0.00
31		3,3'-Dichlorobenzidine	1.000	1.033	-3.3	96	0.00
32		Chrysene	1.000	1.244	-24.4	116	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	1.105	-10.5	109	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	0.916	8.4	89	0.01
35	I	Perylene-d12	5.000	5.000	0.0	92	0.01
36		Benzo(b)fluoranthene	1.000	0.875	12.5	87	0.00
37		Benzo(k)fluoranthene	1.000	0.971	2.9	101	0.00
38	C	Benzo(a)pyrene	1.000	0.905	9.5	94	0.01
39		Dibenzo(a,h)anthracene	1.000	0.859	14.1	88	0.01
40		Benzo(g,h,i)perylene	1.000	0.859	14.1	89	0.01

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

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