Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM012716\
Data File : BM004138.D
Acq On : 27 Jan 2016 17:25
Operator : SJ/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC001

Quant Time: Jan 27 23:59:39 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	74	0.00
2		1,4-Dioxane	1.000	0.984	1.6	76	0.00
3		n-Nitrosodimethylamine	1.000	0.912	8.8	74	0.00
4	S	2-Fluorophenol	1.000	0.914	8.6	75	0.00
5	S	Phenol-d6	1.000	0.928	7.2	77	0.00
6		bis(2-Chloroethyl)ether	1.000	0.926	7.4	76	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	79	0.00
8	S	Nitrobenzene-d5	1.000	0.922	7.8	80	0.00
9		Nitrobenzene	1.000	0.906	9.4	80	0.00
10		Naphthalene	1.000	0.913	8.7	78	0.00
11		Hexachlorobutadiene	1.000	0.903	9.7	78	0.00
12		2-Methylnaphthalene	1.000	0.944	5.6	81	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	86	0.00
14	S	2,4,6-Tribromophenol	1.000	0.934	6.6	90	0.00
15	S	2-Fluorobiphenyl	1.000	0.892	10.8	83	0.00
16		Acenaphthylene	1.000	0.886	11.4	84	0.00
17		Acenaphthene	1.000	0.877	12.3	85	0.00
18		Fluorene	1.000	0.932	6.8	90	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	92	0.00
20		4-Bromophenyl-phenylether	1.000	0.887	11.3	84	0.00
21		Hexachlorobenzene	1.000	0.890	11.0	87	0.00
22		Pentachlorophenol	1.000	0.919	8.1	105	0.00
23		Phenanthrene	1.000	0.923	7.7	92	0.00
24		Anthracene	1.000	0.908	9.2	90	0.00
25		Fluoranthene	1.000	0.940	6.0	94	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	86	0.00
27		Benzidine	1.000	1.008	-0.8	88	0.02
28		Pyrene	1.000	0.955	4.5	90	0.00
29	S	Terphenyl-d14	1.000	0.979	2.1	94	0.00
30		Benzo(a)anthracene	1.000	0.902	9.8	88	0.00
31		3,3'-Dichlorobenzidine	1.000	1.086	-8.6	103	0.00
32		Chrysene	1.000	1.069	-6.9	103	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	1.085	-8.5	109	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	0.954	4.6	94	0.00
35	I	Perylene-d12	5.000	5.000	0.0	94	0.00
36		Benzo(b)fluoranthene	1.000	0.896	10.4	90	0.00
37		Benzo(k)fluoranthene	1.000	0.937	6.3	99	0.00
38	C	Benzo(a)pyrene	1.000	0.910	9.0	96	0.01
39		Dibenzo(a,h)anthracene	1.000	0.903	9.7	94	0.00
40		Benzo(g,h,i)perylene	1.000	0.900	10.0	94	0.00

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

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