Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061516\
Data File : BE091919.D
Acq On : 15 Jun 2016 11:23
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 15 14:21:08 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE061416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 15 12:00:38 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	77	0.00
2		1,4-Dioxane	0.400	0.381	4.8	71	0.00
3		n-Nitrosodimethylamine	0.400	0.332	17.0	64	0.04
4	S	2-Fluorophenol	0.400	0.377	5.8	74	0.02
5	S	Phenol-d6	0.400	0.428	-7.0	76	0.02
6		bis(2-Chloroethyl)ether	0.400	0.347	13.3	67	0.02
7		n-Nitroso-di-n-propylamine	0.400	0.418	-4.5	71	0.02
8	I	Naphthalene-d8	5.000	5.000	0.0	80	0.03
9	S	Nitrobenzene-d5	0.400	0.376	6.0	75	0.02
10		Nitrobenzene	0.400	0.358	10.5	70	0.02
11		bis(2-Chloroethoxy)methane	0.400	0.353	11.8	65	0.03
12		Naphthalene	0.400	0.395	1.3	74	0.00
13		Hexachlorobutadiene	0.400	0.387	3.3	72	0.00
14		2-Methylnaphthalene	0.400	0.383	4.3	74	0.01
15	I	Acenaphthene-d10	5.000	5.000	0.0	81	0.00
16	S	2,4,6-Tribromophenol	0.400	0.350	12.5	74	0.02
17	S	2-Fluorobiphenyl	0.400	0.401	-0.3	75	0.00
18		Acenaphthylene	0.400	0.353	11.8	72	0.03
19		2,6-Dinitrotoluene	0.400	0.468	-17.0	90	0.00
20		Acenaphthene	0.400	0.373	6.8	75	0.00
21		2,4-Dinitrotoluene	0.400	0.379	5.3	73	0.03
22		Fluorene	0.400	0.377	5.8	75	0.02
23		4-Chlorophenyl-phenylether	0.400	0.384	4.0	73	0.02
24	I	Phenanthrene-d10	5.000	5.000	0.0	81	0.00
25		4-Bromophenyl-phenylether	0.400	0.372	7.0	74	0.02
26		Hexachlorobenzene	0.400	0.378	5.5	73	0.00
27		Pentachlorophenol	0.400	0.347	13.3	70	0.02
28		Phenanthrene	0.400	0.387	3.3	75	0.00
29		Anthracene	0.400	0.365	8.8	75	0.00
30		Fluoranthene	0.400	0.376	6.0	75	0.00
31	I	Chrysene-d12	5.000	5.000	0.0	85	0.00
32		Benzidine	0.400	0.382	4.5	98	0.00
33		Pyrene	0.400	0.340	15.0	74	0.00
34	S	Terphenyl-d14	0.400	0.339	15.3	70	0.00
35		Benzo(a)anthracene	0.400	0.383	4.3	83	0.00
36		3,3'-Dichlorobenzidine	0.400	0.324	19.0	83	0.03
37		Chrysene	0.400	0.411	-2.7	78	0.00
38		Bis(2-ethylhexyl)phthalate	0.400	0.395	1.3	85	-0.03
39		Indeno(1,2,3-cd)pyrene	0.400	0.352	12.0	76	0.00
40	I	Perylene-d12	5.000	5.000	0.0	82	0.00
41		Benzo(b)fluoranthene	0.400	0.346	13.5	69	-0.01

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.403	-0.8	83	0.00
43 C	Benzo(a)pyrene	0.400	0.373	6.8	77	0.00
44	Dibenzo(a,h)anthracene	0.400	0.365	8.8	75	-0.02
45	Benzo(g,h,i)perylene	0.400	0.366	8.5	75	0.00

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