Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072716\
Data File : BE092064.D
Acq On : 28 Jul 2016 2:28
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 28 12:41:35 2016
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE072716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 27 17:13:26 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	109	-0.02
2		1,4-Dioxane	0.400	0.373	6.8	102	0.00
3		n-Nitrosodimethylamine	0.400	0.365	8.8	110	0.00
4	S	2-Fluorophenol	0.400	0.416	-4.0	130	-0.02
5	S	Phenol-d6	0.400	0.403	-0.8	124	0.00
6		bis(2-Chloroethyl)ether	0.400	0.378	5.5	114	0.00
7		n-Nitroso-di-n-propylamine	0.400	0.405	-1.3	125	0.00
8	I	Naphthalene-d8	5.000	5.000	0.0	114	0.00
9	S	Nitrobenzene-d5	0.400	0.351	12.3	108	0.00
10		Nitrobenzene	0.400	0.380	5.0	123	0.00
11		bis(2-Chloroethoxy)methane	0.400	0.372	7.0	122	0.00
12		Naphthalene	0.400	0.403	-0.8	116	0.00
13		Hexachlorobutadiene	0.400	0.418	-4.5	122	0.00
14		2-Methylnaphthalene	0.400	0.406	-1.5	124	0.00
15	I	Acenaphthene-d10	5.000	5.000	0.0	115	0.00
16	S	2,4,6-Tribromophenol	0.400	0.483	-20.7	142	0.00
17	S	2-Fluorobiphenyl	0.400	0.407	-1.7	119	0.00
18		Acenaphthylene	0.400	0.406	-1.5	126	0.00
19		2,6-Dinitrotoluene	0.400	0.424	-6.0	115	0.00
20		Acenaphthene	0.400	0.406	-1.5	119	-0.03
21		2,4-Dinitrotoluene	0.400	0.415	-3.7	147	0.00
22		Fluorene	0.400	0.402	-0.5	119	0.00
23		4-Chlorophenyl-phenylether	0.400	0.396	1.0	115	0.00
24	I	Phenanthrene-d10	5.000	5.000	0.0	111	0.00
25		4-Bromophenyl-phenylether	0.400	0.395	1.3	113	0.00
26		Hexachlorobenzene	0.400	0.397	0.8	113	0.00
27		Pentachlorophenol	0.400	0.908	-127.0#	375	0.00
28		Phenanthrene	0.400	0.398	0.5	114	0.00
29		Anthracene	0.400	0.394	1.5	118	0.00
30		Fluoranthene	0.400	0.426	-6.5	135	0.00
31	I	Chrysene-d12	5.000	5.000	0.0	94	0.00
32		Benzidine	0.400	1.495	-273.8#	500	-0.02
33		Pyrene	0.400	0.474	-18.5	122	0.00
34	S	Terphenyl-d14	0.400	0.479	-19.7	127	0.00
35		Benzo(a)anthracene	0.400	0.483	-20.7	112	0.00
36		3,3'-Dichlorobenzidine	0.400	0.666	-66.5#	214	0.00
37		Chrysene	0.400	0.330	17.5	94	0.00
38		Bis(2-ethylhexyl)phthalate	0.400	0.695	-73.7#	177	0.00
39		Indeno(1,2,3-cd)pyrene	0.400	0.528	-32.0#	133	0.00
40	I	Perylene-d12	5.000	5.000	0.0	117	0.00
41		Benzo(b)fluoranthene	0.400	0.412	-3.0	123	0.00

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	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	0.400	0.397	0.8	111	0.00
43 C	Benzo(a)pyrene	0.400	0.414	-3.5	125	0.00
44	Dibenzo(a,h)anthracene	0.400	0.423	-5.7	135	0.00
45	Benzo(g,h,i)perylene	0.400	0.421	-5.2	133	0.00

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