Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE062416\
Data File : BE091956.D
Acq On : 24 Jun 2016 21:07
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 24 23:07:10 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	106	0.02
2		1,4-Dioxane	0.705	0.644	8.7	94	0.00
3		n-Nitrosodimethylamine	0.802	0.778	3.0	102	0.04
4	S	2-Fluorophenol	0.977	1.296	-32.7#	143	0.02
5	S	Phenol-d6	1.267	1.796	-41.8#	157#	0.02
6		bis(2-Chloroethyl)ether	1.431	1.405	1.8	105	0.02
7		n-Nitroso-di-n-propylamine	1.137	1.449	-27.4#	137	0.02
8	I	Naphthalene-d8	1.000	1.000	0.0	120	0.03
9	S	Nitrobenzene-d5	0.345	0.377	-9.3	130	0.02
10		Nitrobenzene	0.343	0.382	-11.4	131	0.02
11		bis(2-Chloroethoxy)methane	0.438	0.499	-13.9	127	0.00
12		Naphthalene	1.092	1.144	-4.8	118	0.00
13		Hexachlorobutadiene	0.177	0.188	-6.2	120	0.00
14		2-Methylnaphthalene	0.741	0.813	-9.7	127	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	115	0.00
16	S	2,4,6-Tribromophenol	0.102	0.140	-37.3#	168#	0.02
17	S	2-Fluorobiphenyl	1.269	1.380	-8.7	116	0.00
18		Acenaphthylene	5.075	5.733	-13.0	131	0.03
19		2,6-Dinitrotoluene	0.811	0.681	16.0	136	0.00
20		Acenaphthene	1.398	1.485	-6.2	121	0.00
21		2,4-Dinitrotoluene	0.758	1.303	-71.9#	211#	0.03
22		Fluorene	1.391	1.478	-6.3	121	0.02
23		4-Chlorophenyl-phenylether	0.685	0.696	-1.6	110	0.00
24	I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
25		4-Bromophenyl-phenylether	0.193	0.228	-18.1	114	0.02
26		Hexachlorobenzene	0.198	0.225	-13.6	107	0.00
27		Pentachlorophenol	0.054	0.070	-29.6#	163#	0.02
28		Phenanthrene	1.250	1.354	-8.3	102	0.00
29		Anthracene	1.106	1.271	-14.9	115	0.00
30		Fluoranthene	5.354	5.697	-6.4	103	0.00
31	I	Chrysene-d12	1.000	1.000	0.0	83	0.00
32		Benzidine	0.170	0.417	-145.3#	286#	0.00
33		Pyrene	3.401	4.074	-19.8	102	0.00
34	S	Terphenyl-d14	0.612	0.663	-8.3	87	0.00
35		Benzo(a)anthracene	0.786	0.984	-25.2#	106	0.00
36		3,3'-Dichlorobenzidine	0.127	0.237	-86.6#	186#	0.03
37		Chrysene	1.147	1.357	-18.3	89	0.00
38		Bis(2-ethylhexyl)phthalate	0.580	1.832	-215.9#	303#	-0.03
39		Indeno(1,2,3-cd)pyrene	3.925	4.647	-18.4	100	0.00
40	I	Perylene-d12	1.000	1.000	0.0	87	0.00
41		Benzo(b)fluoranthene	1.284	1.495	-16.4	98	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	1.175	1.278	-8.8	95	0.00
43 C	Benzo(a)pyrene	1.122	1.328	-18.4	104	0.00
44	Dibenzo(a,h)anthracene	1.092	1.240	-13.6	99	-0.02
45	Benzo(g,h,i)perylene	4.507	5.180	-14.9	99	0.00

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