Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE080116\
Data File : BE092104.D
Acq On : 1 Aug 2016 22:52
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Aug 02 01:28:54 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE080116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 01 17:22:55 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	103	0.00
2		1,4-Dioxane	0.722	0.695	3.7	100	0.00
3		n-Nitrosodimethylamine	0.918	0.806	12.2	96	0.02
4	S	2-Fluorophenol	1.631	1.675	-2.7	102	0.00
5	S	Phenol-d6	1.960	2.050	-4.6	123	0.00
6		bis(2-Chloroethyl)ether	1.545	1.416	8.3	100	0.00
7		n-Nitroso-di-n-propylamine	1.469	1.373	6.5	103	0.00
8	I	Naphthalene-d8	1.000	1.000	0.0	119	0.00
9	S	Nitrobenzene-d5	0.429	0.361	15.9	106	0.00
10		Nitrobenzene	0.416	0.343	17.5	106	0.00
11		bis(2-Chloroethoxy)methane	0.560	0.526	6.1	103	0.00
12		Naphthalene	1.152	1.135	1.5	116	0.00
13		Hexachlorobutadiene	0.188	0.183	2.7	116	0.00
14		2-Methylnaphthalene	0.772	0.770	0.3	121	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	136	0.00
16	S	2,4,6-Tribromophenol	0.363	0.259	28.7#	97	0.00
17	S	2-Fluorobiphenyl	1.936	1.890	2.4	135	0.00
18		Acenaphthylene	11.081	9.318	15.9	118	0.00
19		2,6-Dinitrotoluene	1.299	0.896	31.0#	99	0.00
20		Acenaphthene	1.530	1.560	-2.0	138	0.00
21		2,4-Dinitrotoluene	2.137	1.391	34.9#	104	0.00
22		Fluorene	2.300	1.987	13.6	119	0.00
23		4-Chlorophenyl-phenylether	1.173	1.071	8.7	115	0.00
24	I	Phenanthrene-d10	1.000	1.000	0.0	112	0.00
25		4-Bromophenyl-phenylether	0.168	0.165	1.8	117	0.00
26		Hexachlorobenzene	0.167	0.174	-4.2	119	0.00
27		Pentachlorophenol	0.031	0.012	61.3#	70	0.00
28		Phenanthrene	1.344	0.995	26.0#	84	0.00
29		Anthracene	0.997	0.938	5.9	110	0.00
30		Fluoranthene	3.917	3.813	2.7	108	-0.03
31	I	Chrysene-d12	1.000	1.000	0.0	101	0.00
32		Benzidine	0.390	0.268	31.3#	118	0.00
33		Pyrene	5.683	5.634	0.9	103	0.00
34	S	Terphenyl-d14	0.760	0.611	19.6	72	0.00
35		Benzo(a)anthracene	1.352	1.359	-0.5	112	0.00
36		3,3'-Dichlorobenzidine	0.449	0.400	10.9	95	0.00
37		Chrysene	1.190	1.205	-1.3	96	-0.02
38		Bis(2-ethylhexyl)phthalate	1.208	0.857	29.1#	77	0.00
39		Indeno(1,2,3-cd)pyrene	5.303	4.884	7.9	96	0.00
40	I	Perylene-d12	1.000	1.000	0.0	95	0.00
41		Benzo(b)fluoranthene	1.387	1.378	0.6	100	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	1.419	1.364	3.9	88	0.00
43 C	Benzo(a)pyrene	1.288	1.251	2.9	95	0.00
44	Dibenzo(a,h)anthracene	1.197	1.154	3.6	95	-0.02
45	Benzo(g,h,i)perylene	5.034	5.001	0.7	96	-0.02

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