Evaluate Continuing Calibration Report

Data Path : V:\HPCHEM1\BNA E\DATA\BE070916\
Data File : BE091991.D
Acq On : 9 Jul 2016 10:19
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 11 04:38:01 2016
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE070516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 05 16:14:58 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	105	0.00
2		1,4-Dioxane	0.678	0.653	3.7	97	0.00
3		n-Nitrosodimethylamine	0.768	0.708	7.8	103	0.00
4	S	2-Fluorophenol	1.196	0.976	18.4	91	0.00
5	S	Phenol-d6	1.662	1.360	18.2	95	0.00
6		bis(2-Chloroethyl)ether	1.449	1.397	3.6	109	-0.02
7		n-Nitroso-di-n-propylamine	1.387	1.155	16.7	94	-0.02
8	I	Naphthalene-d8	1.000	1.000	0.0	107	0.00
9	S	Nitrobenzene-d5	0.341	0.298	12.6	95	-0.02
10		Nitrobenzene	0.370	0.321	13.2	99	0.00
11		bis(2-Chloroethoxy)methane	0.418	0.377	9.8	84	-0.03
12		Naphthalene	1.164	1.133	2.7	101	0.00
13		Hexachlorobutadiene	0.177	0.168	5.1	95	0.00
14		2-Methylnaphthalene	0.748	0.700	6.4	98	-0.01
15	I	Acenaphthene-d10	1.000	1.000	0.0	98	0.00
16	S	2,4,6-Tribromophenol	0.143	0.111	22.4	90	0.00
17	S	2-Fluorobiphenyl	1.220	1.241	-1.7	97	0.00
18		Acenaphthylene	6.276	6.257	0.3	101	0.00
19		2,6-Dinitrotoluene	1.093	0.967	11.5	120	0.00
20		Acenaphthene	1.367	1.270	7.1	89	0.00
21		2,4-Dinitrotoluene	1.115	1.249	-12.0	167#	0.00
22		Fluorene	1.442	1.434	0.6	95	0.00
23		4-Chlorophenyl-phenylether	0.719	0.726	-1.0	94	0.00
24	I	Phenanthrene-d10	1.000	1.000	0.0	105	0.00
25		4-Bromophenyl-phenylether	0.217	0.196	9.7	93	0.00
26		Hexachlorobenzene	0.216	0.212	1.9	102	0.00
27		Pentachlorophenol	0.078	0.064	17.9	150	0.00
28		Phenanthrene	1.303	1.274	2.2	99	0.00
29		Anthracene	1.218	1.120	8.0	98	0.00
30		Fluoranthene	5.599	5.233	6.5	99	0.00
31	I	Chrysene-d12	1.000	1.000	0.0	122	0.00
32		Benzidine	0.333	0.254	23.7	110	0.00
33		Pyrene	3.640	3.001	17.6	100	0.00
34	S	Terphenyl-d14	0.590	0.517	12.4	106	0.00
35		Benzo(a)anthracene	0.837	0.771	7.9	113	0.00
36		3,3'-Dichlorobenzidine	0.257	0.156	39.3#	73	0.00
37		Chrysene	1.187	1.184	0.3	122	0.00
38		Bis(2-ethylhexyl)phthalate	0.970	0.588	39.4#	80	-0.03
39		Indeno(1,2,3-cd)pyrene	4.657	3.692	20.7	94	-0.02
40	I	Perylene-d12	1.000	1.000	0.0	107	0.00
41		Benzo(b)fluoranthene	1.381	1.333	3.5	105	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
42	Benzo(k)fluoranthene	1.335	1.179	11.7	96	0.00
43 C	Benzo(a)pyrene	1.304	1.157	11.3	96	0.00
44	Dibenzo(a,h)anthracene	1.227	1.066	13.1	95	0.00
45	Benzo(g,h,i)perylene	5.004	4.301	14.0	93	0.00

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