Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE051017\
Data File : BE092981.D
Acq On : 10 May 2017 14:35
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 11 05:09:42 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE051017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 10 03:43:15 2017
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	99	0.00
2		1,4-Dioxane	0.838	0.769	8.2	99	0.00
3		n-Nitrosodimethylamine	0.692	0.635	8.2	95	0.00
4	S	2-Fluorophenol	1.125	1.053	6.4	99	0.00
5	S	Phenol-d6	1.406	1.312	6.7	101	0.00
6		bis(2-Chloroethyl)ether	1.368	1.263	7.7	96	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
8	S	Nitrobenzene-d5	0.312	0.295	5.4	99	0.00
9		Naphthalene	1.073	1.007	6.2	97	0.00
10		Hexachlorobutadiene	0.153	0.142	7.2	94	0.00
11	SURR	2-Methylnaphthalene-d10	0.560	0.545	2.7	103	-0.02
12		2-Methylnaphthalene	0.616	0.596	3.2	103	0.00
13	I	Acenaphthene-d10	1.000	1.000	0.0	97	0.00
14	S	2,4,6-Tribromophenol	0.104	0.094	9.6	107	-0.01
15	S	2-Fluorobiphenyl	1.498	1.511	-0.9	101	0.00
16		Acenaphthylene	6.761	6.218	8.0	98	0.00
17		Acenaphthene	1.250	1.217	2.6	101	-0.02
18		Fluorene	1.518	1.488	2.0	99	0.00
19	I	Phenanthrene-d10	1.000	1.000	0.0	105	0.00
20		Hexachlorobenzene	0.188	0.158	16.0	95	0.00
21		Pentachlorophenol	0.050	0.046	8.0	112	-0.02
22		Phenanthrene	1.137	1.039	8.6	104	0.00
23		Anthracene	0.901	0.782	13.2	98	0.00
24	SURR	Fluoranthene-d10	1.041	0.925	11.1	103	0.00
25		Fluoranthene	5.367	4.633	13.7	100	0.00
26	I	Chrysene-d12	1.000	1.000	0.0	104	0.00
27		Pyrene	5.349	5.054	5.5	102	0.00
28	S	Terphenyl-d14	0.615	0.587	4.6	103	0.00
29		Benzo(a)anthracene	0.957	0.949	0.8	113	0.00
30		Chrysene	1.279	1.187	7.2	102	0.00
31		Bis(2-ethylhexyl)phthalate	0.429	0.388	9.6	104	0.00
32		Indeno(1,2,3-cd)pyrene	4.419	4.265	3.5	104	-0.02
33	I	Perylene-d12	1.000	1.000	0.0	105	-0.01
34		Benzo(b)fluoranthene	1.226	1.088	11.3	110	0.00
35		Benzo(k)fluoranthene	1.313	1.195	9.0	101	-0.01
36	C	Benzo(a)pyrene	1.145	1.038	9.3	109	0.00
37		Dibenzo(a,h)anthracene	1.133	0.995	12.2	101	0.00
38		Benzo(g,h,i)perylene	4.931	4.276	13.3	100	0.00

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