Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE042418\
Data File : BE096123.D
Acq On : 24 Apr 2018 17:19
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 5 Sample Multiplier: 1

Quant Time: Apr 25 05:24:31 2018
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE042418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 24 17:40:03 2018
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% 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	I	Naphthalene-d8	1.000	1.000	0.0	102	0.00
3		Naphthalene	1.114	1.097	1.5	100	0.00
4	SURR	2-Methylnaphthalene-d10	0.739	0.719	2.7	98	0.00
5		2-Methylnaphthalene	0.746	0.748	-0.3	102	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	101	0.00
7		Acenaphthylene	2.062	2.043	0.9	101	0.00
8	C	Acenaphthene	1.487	1.496	-0.6	101	0.00
9		Fluorene	1.967	2.041	-3.8	121	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	104	0.00
11		Pentachlorophenol	0.096	0.078	18.8	88	0.00
12		Phenanthrene	1.178	1.150	2.4	102	0.00
13		Anthracene	1.182	1.135	4.0	101	0.00
14	SURR	Fluoranthene-d10	1.845	1.783	3.4	101	0.00
15	C	Fluoranthene	1.800	1.763	2.1	103	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	103	0.00
17		Pyrene	0.802	0.752	6.2	97	0.00
18		Benzo(a)anthracene	1.211	1.162	4.0	105	0.00
19		Chrysene	1.113	1.043	6.3	102	0.00
20	I	Perylene-d12	1.000	1.000	0.0	102	0.00
21		Benzo(b)fluoranthene	1.515	1.358	10.4	95	0.00
22		Benzo(k)fluoranthene	1.448	1.306	9.8	97	0.00
23	C	Benzo(a)pyrene	1.316	1.232	6.4	98	0.00
24		Indeno(1,2,3-cd)pyrene	1.773	1.497	15.6	89	0.00
25		Dibenzo(a,h)anthracene	1.485	1.232	17.0	88	0.00
26		Benzo(g,h,i)perylene	1.543	1.255	18.7	87	0.00

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