Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081617\
Data File : BE093846.D
Acq On : 16 Aug 2017 18:48
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTD0.410

Quant Time: Aug 17 01:00:38 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE072417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 17 00:58:19 2017
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	70	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	72	0.00
3		Naphthalene	0.973	0.960	1.3	72	0.00
4	SURR	2-Methylnaphthalene-d10	0.648	0.625	3.5	71	0.00
5		2-Methylnaphthalene	0.710	0.664	6.5	68	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	62	0.00
7		Acenaphthylene	1.673	1.744	-4.2	67	0.00
8	C	Acenaphthene	1.316	1.278	2.9	61	0.00
9		Fluorene	1.664	1.466	11.9	56	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	51	0.00
11		Pentachlorophenol	0.066	0.059	10.6	51	0.00
12		Phenanthrene	1.089	1.030	5.4	50	0.00
13		Anthracene	1.061	1.032	2.7	52	0.00
14	SURR	Fluoranthene-d10	1.263	1.145	9.3	48	0.00
15	C	Fluoranthene	1.362	1.216	10.7	48	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	48	0.00
17		Pyrene	1.168	1.167	0.1	49	0.00
18		Benzo(a)anthracene	1.160	1.118	3.6	48	0.00
19		Chrysene	1.112	1.113	-0.1	49	0.00
20	I	Perylene-d12	1.000	1.000	0.0	48	0.00
21		Benzo(b)fluoranthene	1.237	1.126	9.0	45	0.00
22		Benzo(k)fluoranthene	1.220	1.207	1.1	49	0.00
23	C	Benzo(a)pyrene	1.177	1.151	2.2	49	0.00
24		Indeno(1,2,3-cd)pyrene	1.299	1.200	7.6	46	0.00
25		Dibenzo(a,h)anthracene	1.086	0.995	8.4	46	0.00
26		Benzo(g,h,i)perylene	1.095	0.996	9.0	45	0.00

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