Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081617\
Data File : BE093833.D
Acq On : 16 Aug 2017 10:18
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTD0.409

Quant Time: Aug 17 00:55:24 2017
Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE072417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 16 07:30:49 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	73	0.00
3		Naphthalene	0.973	0.960	1.3	73	0.00
4	SURR	2-Methylnaphthalene-d10	0.648	0.625	3.5	71	0.00
5		2-Methylnaphthalene	0.710	0.662	6.8	69	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	63	0.00
7		Acenaphthylene	1.673	1.684	-0.7	66	0.00
8	C	Acenaphthene	1.316	1.285	2.4	63	0.00
9		Fluorene	1.664	1.476	11.3	58	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	52	0.00
11		Pentachlorophenol	0.066	0.057	13.6	51	0.00
12		Phenanthrene	1.089	1.071	1.7	53	0.00
13		Anthracene	1.061	1.014	4.4	52	0.00
14	SURR	Fluoranthene-d10	1.263	1.123	11.1	48	0.00
15	C	Fluoranthene	1.362	1.194	12.3	48	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	48	0.00
17		Pyrene	1.168	1.149	1.6	49	0.00
18		Benzo(a)anthracene	1.160	1.086	6.4	47	0.00
19		Chrysene	1.112	1.115	-0.3	50	0.00
20	I	Perylene-d12	1.000	1.000	0.0	46	0.00
21		Benzo(b)fluoranthene	1.237	1.148	7.2	45	0.00
22		Benzo(k)fluoranthene	1.220	1.207	1.1	48	0.00
23	C	Benzo(a)pyrene	1.177	1.153	2.0	47	0.00
24		Indeno(1,2,3-cd)pyrene	1.299	1.079	16.9	40	0.00
25		Dibenzo(a,h)anthracene	1.086	0.896	17.5	40	0.00
26		Benzo(g,h,i)perylene	1.095	0.891	18.6	39	0.00

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