Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE080817\
Data File : BE093709.D
Acq On : 8 Aug 2017 10:02
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTD0.444

Quant Time: Aug 09 04:13:35 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 04 16:09:39 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	94	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	91	0.00
3		Naphthalene	0.973	0.982	-0.9	93	0.00
4	SURR	2-Methylnaphthalene-d10	0.648	0.636	1.9	91	0.00
5		2-Methylnaphthalene	0.710	0.690	2.8	90	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	81	0.00
7		Acenaphthylene	1.673	1.659	0.8	83	0.00
8	C	Acenaphthene	1.316	1.300	1.2	82	0.00
9		Fluorene	1.664	1.534	7.8	77	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	74	0.00
11		Pentachlorophenol	0.066	0.044	33.3#	55	0.02
12		Phenanthrene	1.089	1.084	0.5	76	0.00
13		Anthracene	1.061	1.014	4.4	73	0.00
14	SURR	Fluoranthene-d10	1.263	1.122	11.2	68	0.00
15	C	Fluoranthene	1.362	1.214	10.9	69	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	69	0.00
17		Pyrene	1.168	1.138	2.6	69	0.00
18		Benzo(a)anthracene	1.160	1.090	6.0	68	0.00
19		Chrysene	1.112	1.112	0.0	71	0.00
20	I	Perylene-d12	1.000	1.000	0.0	72	0.00
21		Benzo(b)fluoranthene	1.237	1.175	5.0	71	0.00
22		Benzo(k)fluoranthene	1.220	1.157	5.2	70	0.00
23	C	Benzo(a)pyrene	1.177	1.122	4.7	71	0.00
24		Indeno(1,2,3-cd)pyrene	1.299	1.246	4.1	72	0.00
25		Dibenzo(a,h)anthracene	1.086	1.036	4.6	72	0.00
26		Benzo(g,h,i)perylene	1.095	1.018	7.0	69	0.00

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