Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE111417\
Data File : BE094808.D
Acq On : 14 Nov 2017 15:03
Operator : SJ/JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SICV90

Quant Time: Nov 14 16:35:04 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE111417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 14 16:33:16 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	92	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	100	0.00
3		Naphthalene	0.953	0.951	0.2	96	0.01
4	SURR	2-Methylnaphthalene-d10	0.646	0.645	0.2	96	0.01
5		2-Methylnaphthalene	0.663	0.660	0.5	96	0.01
6	I	Acenaphthene-d10	1.000	1.000	0.0	104	0.00
7		Acenaphthylene	1.832	1.799	1.8	99	0.00
8	C	Acenaphthene	1.289	1.291	-0.2	101	0.00
9		Fluorene	1.564	1.577	-0.8	101	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	106	0.00
11		Pentachlorophenol	0.021	0.019	9.5	90	0.02
12		Phenanthrene	1.074	1.070	0.4	104	0.00
13		Anthracene	1.095	1.081	1.3	101	0.00
14	SURR	Fluoranthene-d10	1.374	1.386	-0.9	102	0.00
15	C	Fluoranthene	1.410	1.422	-0.9	102	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	105	0.00
17		Pyrene	1.252	1.244	0.6	103	0.00
18		Benzo(a)anthracene	1.216	1.199	1.4	101	0.00
19		Chrysene	1.214	1.222	-0.7	104	0.00
20	I	Perylene-d12	1.000	1.000	0.0	107	0.00
21		Benzo(b)fluoranthene	1.192	1.196	-0.3	102	0.00
22		Benzo(k)fluoranthene	1.267	1.236	2.4	105	0.00
23	C	Benzo(a)pyrene	1.185	1.169	1.4	103	0.00
24		Indeno(1,2,3-cd)pyrene	1.292	1.272	1.5	102	0.00
25		Dibenzo(a,h)anthracene	1.089	1.070	1.7	101	0.00
26		Benzo(g,h,i)perylene	1.082	1.055	2.5	101	0.00

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