Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM120216\
Data File : BM008179.D
Acq On : 03 Dec 2016 04:46
Operator : UM/SJ
Sample : SSTDCCC0.4EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.464

Quant Time: Dec 03 05:33:31 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM-EPA-SIM-BM112116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Dec 03 04:54:05 2016
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	85	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	88	0.00
3		Naphthalene	1.119	1.063	5.0	83	0.00
4	SURR	2-Methylnaphthalene-d10	0.566	0.535	5.5	83	0.00
5		2-Methylnaphthalene	0.721	0.675	6.4	82	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	78	0.00
7		Acenaphthylene	1.810	1.864	-3.0	81	0.00
8	C	Acenaphthene	1.377	1.336	3.0	78	0.00
9		Fluorene	1.583	1.467	7.3	74	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	71	0.00
11		Pentachlorophenol	0.140	0.116	17.1	61	-0.02
12		Phenanthrene	1.246	1.185	4.9	68	0.00
13		Anthracene	1.170	1.269	-8.5	78	0.00
14	SURR	Fluoranthene-d10	1.132	1.024	9.5	65	0.00
15	C	Fluoranthene	1.608	1.395	13.2	64	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	57	0.00
17		Pyrene	1.389	1.588	-14.3	65	-0.01
18		Benzo(a)anthracene	1.228	1.317	-7.2	63	-0.01
19		Chrysene	1.527	1.413	7.5	54	0.00
20	I	Perylene-d12	1.000	1.000	0.0	63	0.00
21		Benzo(b)fluoranthene	1.596	1.433	10.2	58	0.00
22		Benzo(k)fluoranthene	1.647	1.491	9.5	59	0.00
23	C	Benzo(a)pyrene	1.495	1.387	7.2	60	0.00
24		Indeno(1,2,3-cd)pyrene	1.848	1.780	3.7	61	0.00
25		Dibenzo(a,h)anthracene	1.481	1.415	4.5	61	0.00
26		Benzo(g,h,i)perylene	1.640	1.564	4.6	60	0.00

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