Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM120916\
Data File : BM008208.D
Acq On : 09 Dec 2016 16:51
Operator : UM/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD0.437

Quant Time: Dec 09 17:59:56 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM112116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 09 17:57:55 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	93	0.02
2	I	Naphthalene-d8	1.000	1.000	0.0	104	0.01
3		Naphthalene	1.119	0.954	14.7	88	0.01
4	SURR	2-Methylnaphthalene-d10	0.566	0.497	12.2	91	0.02
5		2-Methylnaphthalene	0.721	0.622	13.7	90	0.01
6	I	Acenaphthene-d10	1.000	1.000	0.0	105	0.00
7		Acenaphthylene	1.810	1.556	14.0	91	0.02
8	C	Acenaphthene	1.377	1.194	13.3	93	0.02
9		Fluorene	1.583	1.307	17.4	88	0.02
10	I	Phenanthrene-d10	1.000	1.000	0.0	93	0.00
11		Pentachlorophenol	0.140	0.101	27.9#	69	0.02
12		Phenanthrene	1.246	1.043	16.3	79	0.02
13		Anthracene	1.170	1.002	14.4	80	0.02
14	SURR	Fluoranthene-d10	1.132	1.008	11.0	84	0.00
15	C	Fluoranthene	1.608	1.302	19.0	78	0.01
16	I	Chrysene-d12	1.000	1.000	0.0	63	0.00
17		Pyrene	1.389	1.575	-13.4	72	0.01
18		Benzo(a)anthracene	1.228	1.216	1.0	64	0.01
19		Chrysene	1.527	1.181	22.7#	50	0.01
20	I	Perylene-d12	1.000	1.000	0.0	67	0.01
21		Benzo(b)fluoranthene	1.596	1.349	15.5	58	0.02
22		Benzo(k)fluoranthene	1.647	1.264	23.3#	53	0.02
23	C	Benzo(a)pyrene	1.495	1.251	16.3	57	0.01
24		Indeno(1,2,3-cd)pyrene	1.848	1.583	14.3	57	0.02
25		Dibenzo(a,h)anthracene	1.481	1.276	13.8	58	0.03
26		Benzo(g,h,i)perylene	1.640	1.385	15.5	56	0.02

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