Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM021715\
Data File : BM000537.D
Acq On : 17 Feb 2015 12:13
Operator : TP/IZ
Sample : SSTD0.415
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC0.4

Quant Time: Feb 18 03:50:53 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM01.2-EPA-SIM-BM021215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Feb 12 16:45:55 2015
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	100	-0.02
2	I	Naphthalene-d8	1.000	1.000	0.0	100	-0.02
3		Naphthalene	0.938	0.932	0.6	98	-0.02
4	SURR	2-Methylnaphthalene-d10	0.510	0.492	3.5	96	-0.02
5		2-Methylnaphthalene	0.634	0.613	3.3	96	-0.02
6	I	Acenaphthene-d10	1.000	1.000	0.0	94	-0.02
7		Acenaphthylene	1.505	1.458	3.1	94	-0.02
8	C	Acenaphthene	1.155	1.133	1.9	92	-0.03
9		Fluorene	1.396	1.328	4.9	91	-0.02
10	I	Phenanthrene-d10	1.000	1.000	0.0	89	-0.02
11		Pentachlorophenol	0.051	0.032	37.3#	61	-0.03
12		Phenanthrene	1.086	1.064	2.0	89	-0.02
13		Anthracene	0.977	0.951	2.7	88	-0.02
14	SURR	Fluoranthene-d10	0.841	0.828	1.5	86	-0.01
15	C	Fluoranthene	1.211	1.169	3.5	84	-0.02
16	I	Chrysene-d12	1.000	1.000	0.0	82	-0.02
17		Pyrene	1.439	1.424	1.0	85	-0.02
18		Benzo(a)anthracene	1.128	1.099	2.6	81	-0.02
19		Chrysene	1.300	1.282	1.4	79	-0.02
20	I	Perylene-d12	1.000	1.000	0.0	79	-0.02
21		Benzo(b)fluoranthene	1.352	1.382	-2.2	80	-0.02
22		Benzo(k)fluoranthene	1.345	1.291	4.0	78	-0.02
23	C	Benzo(a)pyrene	1.264	1.246	1.4	79	-0.03
24		Indeno(1,2,3-cd)pyrene	1.410	1.397	0.9	78	-0.03
25		Dibenzo(a,h)anthracene	1.114	1.119	-0.4	80	-0.04
26		Benzo(g,h,i)perylene	1.244	1.232	1.0	78	-0.04

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