Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072417\
Data File : BE093535.D
Acq On : 24 Jul 2017 13:11
Operator : SJ/JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SICV35

Quant Time: Jul 24 14:28:57 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 24 13:14:05 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	97	0.00
2	I	Naphthalene-d8	1.000	1.000	0.0	96	0.00
3		Naphthalene	0.973	0.987	-1.4	99	0.00
4	SURR	2-Methylnaphthalene-d10	0.648	0.664	-2.5	100	0.00
5		2-Methylnaphthalene	0.710	0.725	-2.1	99	0.00
6	I	Acenaphthene-d10	1.000	1.000	0.0	97	0.00
7		Acenaphthylene	1.673	1.658	0.9	100	0.00
8	C	Acenaphthene	1.316	1.308	0.6	99	0.00
9		Fluorene	1.664	1.646	1.1	99	0.00
10	I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
11		Pentachlorophenol	0.066	0.056	15.2	92	0.00
12		Phenanthrene	1.089	1.078	1.0	100	0.00
13		Anthracene	1.061	1.012	4.6	97	0.00
14	SURR	Fluoranthene-d10	1.263	1.210	4.2	97	0.00
15	C	Fluoranthene	1.362	1.295	4.9	97	0.00
16	I	Chrysene-d12	1.000	1.000	0.0	96	0.00
17		Pyrene	1.168	1.166	0.2	98	0.00
18		Benzo(a)anthracene	1.160	1.131	2.5	97	0.00
19		Chrysene	1.112	1.112	0.0	98	0.00
20	I	Perylene-d12	1.000	1.000	0.0	95	0.01
21		Benzo(b)fluoranthene	1.237	1.246	-0.7	99	0.00
22		Benzo(k)fluoranthene	1.220	1.208	1.0	97	0.00
23	C	Benzo(a)pyrene	1.177	1.178	-0.1	98	0.01
24		Indeno(1,2,3-cd)pyrene	1.299	1.288	0.8	98	0.01
25		Dibenzo(a,h)anthracene	1.086	1.087	-0.1	100	0.00
26		Benzo(g,h,i)perylene	1.095	1.085	0.9	98	0.00

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