Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE090415\
Data File : BE090728.D
Acq On : 4 Sep 2015 14:00
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 04 14:33:07 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 02 11:40:50 2015
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	5.000	5.000	0.0	97	0.00
2		1,4-Dioxane	1.000	0.974	2.6	89	0.00
3		n-Nitrosodimethylamine	1.000	0.940	6.0	88	0.00
4	S	2-Fluorophenol	1.000	1.201	-20.1	115	0.00
5	S	Phenol-d6	1.000	1.161	-16.1	118	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	106	0.00
7	S	Nitrobenzene-d5	1.000	1.016	-1.6	108	0.00
8		Nitrobenzene	1.000	0.977	2.3	109	0.00
9		Naphthalene	1.000	1.002	-0.2	107	0.00
10		Hexachlorobutadiene	1.000	0.952	4.8	98	0.00
11		2-Methylnaphthalene	1.000	1.148	-14.8	126	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	118	0.00
13	S	2,4,6-Tribromophenol	1.000	1.311	-31.1#	170	0.00
14	S	2-Fluorobiphenyl	1.000	1.009	-0.9	121	0.00
15		Acenaphthylene	1.000	1.074	-7.4	132	0.00
16		Acenaphthene	1.000	1.014	-1.4	119	0.00
17		Fluorene	1.000	1.006	-0.6	120	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	111	0.00
19		4-Bromophenyl-phenylether	1.000	1.107	-10.7	128	0.00
20		Hexachlorobenzene	1.000	1.028	-2.8	111	0.00
21		Pentachlorophenol	1.000	1.379	-37.9#	159	0.00
22		Phenanthrene	1.000	1.054	-5.4	117	-0.02
23		Anthracene	1.000	1.120	-12.0	129	0.00
24		Fluoranthene	1.000	1.095	-9.5	131	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	107	0.00
26		Pyrene	1.000	1.200	-20.0	136	0.00
27	S	Terphenyl-d14	1.000	1.160	-16.0	130	0.00
28		Benzo(a)anthracene	1.000	1.095	-9.5	119	0.00
29		Chrysene	1.000	1.105	-10.5	114	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.500	-50.0#	191	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.144	-14.4	133	0.00
32	I	Perylene-d12	5.000	5.000	0.0	116	0.00
33		Benzo(b)fluoranthene	1.000	1.046	-4.6	124	0.00
34		Benzo(k)fluoranthene	1.000	1.089	-8.9	128	0.00
35	C	Benzo(a)pyrene	1.000	1.147	-14.7	142	0.00
36		Dibenzo(a,h)anthracene	1.000	1.063	-6.3	133	0.00
37		Benzo(g,h,i)perylene	1.000	1.063	-6.3	130	-0.01

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