Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE090815\
Data File : BE090755.D
Acq On : 8 Sep 2015 19:15
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 08 23:08:18 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 08 23:05:35 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	133	0.00
2		1,4-Dioxane	1.000	0.935	6.5	118	0.00
3		n-Nitrosodimethylamine	1.000	0.956	4.4	123	0.00
4	S	2-Fluorophenol	1.000	1.039	-3.9	136	0.00
5	S	Phenol-d6	1.000	1.136	-13.6	159	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	148	0.00
7	S	Nitrobenzene-d5	1.000	1.027	-2.7	152	0.00
8		Nitrobenzene	1.000	1.014	-1.4	158	0.00
9		Naphthalene	1.000	1.009	-0.9	150	0.00
10		Hexachlorobutadiene	1.000	0.960	4.0	137	0.00
11		2-Methylnaphthalene	1.000	1.125	-12.5	171	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	172	0.00
13	S	2,4,6-Tribromophenol	1.000	1.102	-10.2	202	0.00
14	S	2-Fluorobiphenyl	1.000	0.968	3.2	169	0.00
15		Acenaphthylene	1.000	1.032	-3.2	185	0.00
16		Acenaphthene	1.000	0.998	0.2	170	0.00
17		Fluorene	1.000	1.012	-1.2	176	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	164	0.00
19		4-Bromophenyl-phenylether	1.000	1.033	-3.3	176	-0.02
20		Hexachlorobenzene	1.000	1.000	0.0	160	0.00
21		Pentachlorophenol	1.000	1.158	-15.8	178	0.00
22		Phenanthrene	1.000	1.029	-2.9	169	-0.02
23		Anthracene	1.000	1.113	-11.3	190	0.00
24		Fluoranthene	1.000	1.031	-3.1	182	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	157	0.00
26		Pyrene	1.000	1.167	-16.7	193	0.00
27	S	Terphenyl-d14	1.000	1.118	-11.8	184	0.00
28		Benzo(a)anthracene	1.000	1.086	-8.6	173	0.00
29		Chrysene	1.000	1.101	-10.1	166	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.107	-10.7	198	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.134	-13.4	193	0.00
32	I	Perylene-d12	5.000	5.000	0.0	167	0.00
33		Benzo(b)fluoranthene	1.000	1.067	-6.7	183	0.00
34		Benzo(k)fluoranthene	1.000	1.037	-3.7	175	0.00
35	C	Benzo(a)pyrene	1.000	1.082	-8.2	193	0.00
36		Dibenzo(a,h)anthracene	1.000	1.063	-6.3	191	0.00
37		Benzo(g,h,i)perylene	1.000	1.023	-2.3	180	-0.01

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