Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE082015\
Data File : BE090571.D
Acq On : 20 Aug 2015 12:05
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 21 01:28:33 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE081415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 14 16:52:42 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	92	0.00
2		1,4-Dioxane	1.000	1.244	-24.4	98	0.00
3		n-Nitrosodimethylamine	1.000	1.076	-7.6	101	0.00
4	S	2-Fluorophenol	1.000	0.999	0.1	92	0.00
5	S	Phenol-d6	1.000	0.927	7.3	89	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	86	0.00
7	S	Nitrobenzene-d5	1.000	1.085	-8.5	99	0.00
8		Nitrobenzene	1.000	1.071	-7.1	99	0.00
9		Naphthalene	1.000	0.974	2.6	85	0.00
10		Hexachlorobutadiene	1.000	0.987	1.3	85	0.00
11		2-Methylnaphthalene	1.000	1.045	-4.5	92	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	98	0.00
13	S	2,4,6-Tribromophenol	1.000	1.425	-42.5#	154	-0.02
14	S	2-Fluorobiphenyl	1.000	0.973	2.7	95	0.00
15		Acenaphthylene	1.000	1.005	-0.5	100	0.00
16		Acenaphthene	1.000	0.976	2.4	97	0.00
17		Fluorene	1.000	1.063	-6.3	105	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	109	0.00
19		4-Bromophenyl-phenylether	1.000	1.018	-1.8	114	0.00
20		Hexachlorobenzene	1.000	0.969	3.1	106	0.00
21		Pentachlorophenol	1.000	1.504	-50.4#	254	0.00
22		Phenanthrene	1.000	0.962	3.8	106	0.00
23		Anthracene	1.000	1.062	-6.2	119	0.00
24		Fluoranthene	1.000	1.149	-14.9	128	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	132	0.00
26		Pyrene	1.000	0.967	3.3	131	0.00
27	S	Terphenyl-d14	1.000	1.011	-1.1	136	0.00
28		Benzo(a)anthracene	1.000	1.023	-2.3	142	0.00
29		Chrysene	1.000	1.007	-0.7	134	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.338	-33.8#	215	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.043	-4.3	149	0.00
32	I	Perylene-d12	5.000	5.000	0.0	139	0.00
33		Benzo(b)fluoranthene	1.000	1.031	-3.1	147	0.00
34		Benzo(k)fluoranthene	1.000	1.061	-6.1	146	0.00
35	C	Benzo(a)pyrene	1.000	1.054	-5.4	151	0.00
36		Dibenzo(a,h)anthracene	1.000	1.042	-4.2	152	0.00
37		Benzo(g,h,i)perylene	1.000	1.040	-4.0	147	0.00

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