Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081315\
Data File : BE090499.D
Acq On : 13 Aug 2015 13:37
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 13 15:11:26 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 06 12:28:08 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	130	-0.01
2		1,4-Dioxane	1.000	1.092	-9.2	125	0.00
3		n-Nitrosodimethylamine	1.000	1.119	-11.9	146	0.00
4	S	2-Fluorophenol	1.000	1.994	-99.4#	278	0.00
5	S	Phenol-d6	1.000	1.561	-56.1#	249	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	145	0.00
7	S	Nitrobenzene-d5	1.000	1.152	-15.2	186	-0.01
8		Nitrobenzene	1.000	1.137	-13.7	196	-0.01
9		Naphthalene	1.000	1.025	-2.5	150	0.00
10		Hexachlorobutadiene	1.000	0.934	6.6	125	0.00
11		2-Methylnaphthalene	1.000	1.159	-15.9	176	-0.01
12	I	Acenaphthene-d10	5.000	5.000	0.0	144	0.00
13	S	2,4,6-Tribromophenol	1.000	2.140	-114.0#	402	-0.02
14	S	2-Fluorobiphenyl	1.000	1.093	-9.3	155	0.00
15		Acenaphthylene	1.000	1.126	-12.6	163	0.00
16		Acenaphthene	1.000	1.090	-9.0	156	0.00
17		Fluorene	1.000	1.029	-2.9	146	-0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	134	0.00
19		4-Bromophenyl-phenylether	1.000	1.060	-6.0	146	0.00
20		Hexachlorobenzene	1.000	0.900	10.0	120	0.00
21		Pentachlorophenol	1.000	2.019	-101.9#	401	0.00
22		Phenanthrene	1.000	1.041	-4.1	141	0.00
23		Anthracene	1.000	1.176	-17.6	162	-0.02
24		Fluoranthene	1.000	1.153	-15.3	161	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	118	0.00
26		Pyrene	1.000	1.337	-33.7#	164	0.00
27	S	Terphenyl-d14	1.000	1.255	-25.5#	150	0.00
28		Benzo(a)anthracene	1.000	1.170	-17.0	154	0.00
29		Chrysene	1.000	0.998	0.2	115	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	2.379	-137.9#	395	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.259	-25.9#	195	0.00
32	I	Perylene-d12	5.000	5.000	0.0	128	0.00
33		Benzo(b)fluoranthene	1.000	1.216	-21.6	178	-0.01
34		Benzo(k)fluoranthene	1.000	0.969	3.1	121	0.00
35	C	Benzo(a)pyrene	1.000	1.220	-22.0#	178	0.00
36		Dibenzo(a,h)anthracene	1.000	1.253	-25.3#	208	-0.01
37		Benzo(g,h,i)perylene	1.000	1.291	-29.1#	171	-0.01

(#) = Out of Range SPCC's out = 0 CCC's out = 1