Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081115\
Data File : BE090429.D
Acq On : 11 Aug 2015 15:40
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 12 04:18:51 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	120	0.00
2		1,4-Dioxane	1.000	1.092	-9.2	116	0.00
3		n-Nitrosodimethylamine	1.000	1.037	-3.7	125	0.00
4	S	2-Fluorophenol	1.000	1.395	-39.5#	180	0.00
5	S	Phenol-d6	1.000	1.213	-21.3	169	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	128	0.00
7	S	Nitrobenzene-d5	1.000	1.024	-2.4	142	0.00
8		Nitrobenzene	1.000	0.991	0.9	145	0.00
9		Naphthalene	1.000	1.013	-1.3	130	0.00
10		Hexachlorobutadiene	1.000	1.074	-7.4	125	0.00
11		2-Methylnaphthalene	1.000	1.118	-11.8	149	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	140	0.00
13	S	2,4,6-Tribromophenol	1.000	1.464	-46.4#	245	0.00
14	S	2-Fluorobiphenyl	1.000	1.045	-4.5	144	0.00
15		Acenaphthylene	1.000	1.057	-5.7	149	0.00
16		Acenaphthene	1.000	1.001	-0.1	140	0.00
17		Fluorene	1.000	1.018	-1.8	141	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	134	0.00
19		4-Bromophenyl-phenylether	1.000	1.001	-0.1	137	0.00
20		Hexachlorobenzene	1.000	0.932	6.8	125	0.00
21		Pentachlorophenol	1.000	1.082	-8.2	176	0.00
22		Phenanthrene	1.000	0.996	0.4	135	0.00
23		Anthracene	1.000	1.096	-9.6	151	-0.02
24		Fluoranthene	1.000	1.075	-7.5	150	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	131	0.00
26		Pyrene	1.000	1.108	-10.8	150	0.00
27	S	Terphenyl-d14	1.000	1.125	-12.5	148	0.00
28		Benzo(a)anthracene	1.000	1.055	-5.5	153	0.00
29		Chrysene	1.000	1.020	-2.0	130	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.177	-17.7	182	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.054	-5.4	172	0.00
32	I	Perylene-d12	5.000	5.000	0.0	136	0.00
33		Benzo(b)fluoranthene	1.000	1.061	-6.1	160	0.00
34		Benzo(k)fluoranthene	1.000	0.969	3.1	128	0.00
35	C	Benzo(a)pyrene	1.000	1.047	-4.7	156	0.00
36		Dibenzo(a,h)anthracene	1.000	1.067	-6.7	178	0.00
37		Benzo(g,h,i)perylene	1.000	1.073	-7.3	151	0.00

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