Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE081115\
Data File : BE090460.D
Acq On : 12 Aug 2015 10:37
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 12 11:25:24 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	121	0.00
2		1,4-Dioxane	1.000	1.031	-3.1	110	0.00
3		n-Nitrosodimethylamine	1.000	1.070	-7.0	130	0.00
4	S	2-Fluorophenol	1.000	1.428	-42.8#	185	0.00
5	S	Phenol-d6	1.000	1.241	-24.1	175	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	128	0.00
7	S	Nitrobenzene-d5	1.000	1.047	-4.7	146	0.00
8		Nitrobenzene	1.000	1.025	-2.5	151	0.00
9		Naphthalene	1.000	1.022	-2.2	131	0.00
10		Hexachlorobutadiene	1.000	1.090	-9.0	127	0.00
11		2-Methylnaphthalene	1.000	1.082	-8.2	144	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	136	0.00
13	S	2,4,6-Tribromophenol	1.000	1.489	-48.9#	244	0.00
14	S	2-Fluorobiphenyl	1.000	1.035	-3.5	139	0.00
15		Acenaphthylene	1.000	1.038	-3.8	143	0.00
16		Acenaphthene	1.000	1.005	-0.5	137	0.00
17		Fluorene	1.000	1.008	-0.8	136	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	132	0.00
19		4-Bromophenyl-phenylether	1.000	1.003	-0.3	136	0.00
20		Hexachlorobenzene	1.000	0.927	7.3	122	0.00
21		Pentachlorophenol	1.000	1.294	-29.4#	219	0.00
22		Phenanthrene	1.000	1.006	-0.6	134	0.00
23		Anthracene	1.000	1.103	-10.3	150	-0.02
24		Fluoranthene	1.000	1.057	-5.7	145	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	129	0.00
26		Pyrene	1.000	1.104	-10.4	148	0.00
27	S	Terphenyl-d14	1.000	1.109	-10.9	145	-0.01
28		Benzo(a)anthracene	1.000	1.072	-7.2	154	0.00
29		Chrysene	1.000	0.983	1.7	124	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.211	-21.1	186	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.050	-5.0	169	0.00
32	I	Perylene-d12	5.000	5.000	0.0	135	0.00
33		Benzo(b)fluoranthene	1.000	1.083	-8.3	163	0.00
34		Benzo(k)fluoranthene	1.000	0.896	10.4	117	0.00
35	C	Benzo(a)pyrene	1.000	1.052	-5.2	156	0.00
36		Dibenzo(a,h)anthracene	1.000	1.065	-6.5	176	-0.01
37		Benzo(g,h,i)perylene	1.000	1.051	-5.1	146	0.00

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