Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE082815\
Data File : BE090667.D
Acq On : 28 Aug 2015 21:54
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 31 04:53:35 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	101	0.00
2		1,4-Dioxane	1.000	1.302	-30.2#	112	0.00
3		n-Nitrosodimethylamine	1.000	1.002	-0.2	103	0.00
4	S	2-Fluorophenol	1.000	0.710	29.0#	72	0.00
5	S	Phenol-d6	1.000	0.668	33.2#	70	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	93	0.00
7	S	Nitrobenzene-d5	1.000	0.843	15.7	82	0.00
8		Nitrobenzene	1.000	0.833	16.7	83	0.00
9		Naphthalene	1.000	0.980	2.0	92	0.00
10		Hexachlorobutadiene	1.000	1.101	-10.1	102	0.00
11		2-Methylnaphthalene	1.000	0.896	10.4	85	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	102	0.00
13	S	2,4,6-Tribromophenol	1.000	0.754	24.6	85	0.00
14	S	2-Fluorobiphenyl	1.000	0.880	12.0	90	0.00
15		Acenaphthylene	1.000	0.894	10.6	93	0.00
16		Acenaphthene	1.000	0.974	2.6	100	0.00
17		Fluorene	1.000	1.038	-3.8	106	0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	118	0.00
19		4-Bromophenyl-phenylether	1.000	0.888	11.2	108	0.00
20		Hexachlorobenzene	1.000	1.004	-0.4	120	-0.02
21		Pentachlorophenol	1.000	0.745	25.5#	102	0.00
22		Phenanthrene	1.000	0.945	5.5	114	0.00
23		Anthracene	1.000	0.912	8.8	112	0.00
24		Fluoranthene	1.000	0.911	8.9	111	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	127	0.00
26		Pyrene	1.000	0.877	12.3	114	0.00
27	S	Terphenyl-d14	1.000	0.855	14.5	111	0.00
28		Benzo(a)anthracene	1.000	0.858	14.2	115	0.00
29		Chrysene	1.000	1.017	-1.7	131	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.549	45.1#	71	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.812	18.8	112	0.00
32	I	Perylene-d12	5.000	5.000	0.0	120	0.00
33		Benzo(b)fluoranthene	1.000	0.922	7.8	113	0.00
34		Benzo(k)fluoranthene	1.000	1.026	-2.6	122	0.00
35	C	Benzo(a)pyrene	1.000	0.900	10.0	111	0.00
36		Dibenzo(a,h)anthracene	1.000	0.837	16.3	106	0.00
37		Benzo(g,h,i)perylene	1.000	0.874	12.6	106	0.00

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