Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE082815\
Data File : BE090656.D
Acq On : 28 Aug 2015 11:51
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 31 04:57:43 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	103	0.00
2		1,4-Dioxane	1.000	1.275	-27.5#	112	0.00
3		n-Nitrosodimethylamine	1.000	0.988	1.2	103	0.00
4	S	2-Fluorophenol	1.000	0.618	38.2#	63	0.00
5	S	Phenol-d6	1.000	0.627	37.3#	67	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	94	0.00
7	S	Nitrobenzene-d5	1.000	0.886	11.4	88	0.00
8		Nitrobenzene	1.000	0.843	15.7	85	0.00
9		Naphthalene	1.000	0.989	1.1	94	0.00
10		Hexachlorobutadiene	1.000	1.084	-8.4	102	0.00
11		2-Methylnaphthalene	1.000	0.891	10.9	86	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	102	0.00
13	S	2,4,6-Tribromophenol	1.000	0.629	37.1#	71	0.00
14	S	2-Fluorobiphenyl	1.000	0.881	11.9	90	0.00
15		Acenaphthylene	1.000	0.882	11.8	92	0.00
16		Acenaphthene	1.000	0.982	1.8	102	0.00
17		Fluorene	1.000	1.046	-4.6	108	0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	120	0.00
19		4-Bromophenyl-phenylether	1.000	0.896	10.4	111	0.00
20		Hexachlorobenzene	1.000	1.007	-0.7	121	-0.02
21		Pentachlorophenol	1.000	0.646	35.4#	80	0.00
22		Phenanthrene	1.000	0.958	4.2	117	0.00
23		Anthracene	1.000	0.888	11.2	110	0.00
24		Fluoranthene	1.000	0.914	8.6	113	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	129	0.00
26		Pyrene	1.000	0.876	12.4	116	0.00
27	S	Terphenyl-d14	1.000	0.862	13.8	113	0.00
28		Benzo(a)anthracene	1.000	0.892	10.8	121	0.00
29		Chrysene	1.000	0.992	0.8	130	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.495	50.5#	62	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.794	20.6	111	0.00
32	I	Perylene-d12	5.000	5.000	0.0	120	0.00
33		Benzo(b)fluoranthene	1.000	0.937	6.3	116	0.00
34		Benzo(k)fluoranthene	1.000	0.998	0.2	119	0.00
35	C	Benzo(a)pyrene	1.000	0.888	11.2	110	0.00
36		Dibenzo(a,h)anthracene	1.000	0.824	17.6	105	0.00
37		Benzo(g,h,i)perylene	1.000	0.856	14.4	105	0.00

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