Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE082515\
Data File : BE090610.D
Acq On : 25 Aug 2015 12:23
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 26 03:25:40 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	104	0.00
2		1,4-Dioxane	1.000	1.350	-35.0#	120	0.00
3		n-Nitrosodimethylamine	1.000	1.022	-2.2	108	0.00
4	S	2-Fluorophenol	1.000	0.764	23.6	79	0.00
5	S	Phenol-d6	1.000	0.724	27.6#	79	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	96	0.00
7	S	Nitrobenzene-d5	1.000	0.929	7.1	94	0.00
8		Nitrobenzene	1.000	0.900	10.0	92	0.00
9		Naphthalene	1.000	0.982	1.8	95	0.00
10		Hexachlorobutadiene	1.000	1.023	-2.3	98	0.00
11		2-Methylnaphthalene	1.000	0.950	5.0	93	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	104	0.00
13	S	2,4,6-Tribromophenol	1.000	0.832	16.8	95	0.00
14	S	2-Fluorobiphenyl	1.000	0.941	5.9	98	0.00
15		Acenaphthylene	1.000	0.940	6.0	99	0.00
16		Acenaphthene	1.000	0.965	3.5	101	0.00
17		Fluorene	1.000	1.050	-5.0	110	0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	120	0.00
19		4-Bromophenyl-phenylether	1.000	0.884	11.6	109	0.00
20		Hexachlorobenzene	1.000	0.977	2.3	118	-0.02
21		Pentachlorophenol	1.000	0.802	19.8	117	0.00
22		Phenanthrene	1.000	0.950	5.0	116	0.00
23		Anthracene	1.000	0.946	5.4	118	0.00
24		Fluoranthene	1.000	1.018	-1.8	126	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	141	0.00
26		Pyrene	1.000	0.900	10.0	130	0.00
27	S	Terphenyl-d14	1.000	0.894	10.6	128	0.00
28		Benzo(a)anthracene	1.000	0.938	6.2	139	0.00
29		Chrysene	1.000	0.994	0.6	142	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.696	30.4#	107	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.894	10.6	137	0.00
32	I	Perylene-d12	5.000	5.000	0.0	140	0.00
33		Benzo(b)fluoranthene	1.000	0.988	1.2	142	0.00
34		Benzo(k)fluoranthene	1.000	0.995	0.5	138	0.00
35	C	Benzo(a)pyrene	1.000	0.949	5.1	138	0.00
36		Dibenzo(a,h)anthracene	1.000	0.900	10.0	133	0.00
37		Benzo(g,h,i)perylene	1.000	0.928	7.2	133	0.00

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