Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE063015\
Data File : BE090027.D
Acq On : 1 Jul 2015 1:08
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 01 05:20:27 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 30 07:01:01 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	90	0.00
2		1,4-Dioxane	1.000	0.849	15.1	80	-0.01
3		n-Nitrosodimethylamine	1.000	0.834	16.6	78	-0.01
4	S	2-Fluorophenol	1.000	1.018	-1.8	92	0.00
5	S	Phenol-d6	1.000	1.007	-0.7	92	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	92	-0.01
7	S	Nitrobenzene-d5	1.000	0.900	10.0	84	0.00
8		Nitrobenzene	1.000	0.884	11.6	84	0.00
9		Naphthalene	1.000	0.922	7.8	86	0.00
10		Hexachlorobutadiene	1.000	0.937	6.3	87	0.00
11		2-Methylnaphthalene	1.000	0.892	10.8	83	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	85	0.00
13	S	2,4,6-Tribromophenol	1.000	1.051	-5.1	92	0.00
14	S	2-Fluorobiphenyl	1.000	0.922	7.8	81	0.00
15		Acenaphthylene	1.000	0.946	5.4	83	0.00
16		Acenaphthene	1.000	0.945	5.5	82	-0.01
17		Fluorene	1.000	0.896	10.4	78	-0.02
18	I	Phenanthrene-d10	5.000	5.000	0.0	77	-0.02
19		4-Bromophenyl-phenylether	1.000	0.895	10.5	71	0.00
20		Hexachlorobenzene	1.000	0.902	9.8	72	-0.02
21		Pentachlorophenol	1.000	0.940	6.0	76	0.00
22		Phenanthrene	1.000	0.895	10.5	71	0.00
23		Anthracene	1.000	0.913	8.7	72	0.00
24		Fluoranthene	1.000	0.952	4.8	75	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	81	-0.01
26		Pyrene	1.000	0.919	8.1	76	0.00
27	S	Terphenyl-d14	1.000	0.890	11.0	72	0.00
28		Benzo(a)anthracene	1.000	0.949	5.1	78	0.00
29		Chrysene	1.000	0.889	11.1	73	-0.01
30		Bis(2-ethylhexyl)phthalate	1.000	1.250	-25.0	128	-0.01
31		Indeno(1,2,3-cd)pyrene	1.000	1.083	-8.3	89	-0.03
32	I	Perylene-d12	5.000	5.000	0.0	83	-0.02
33		Benzo(b)fluoranthene	1.000	1.051	-5.1	88	-0.01
34		Benzo(k)fluoranthene	1.000	0.980	2.0	81	-0.01
35	C	Benzo(a)pyrene	1.000	1.056	-5.6	88	-0.02
36		Dibenzo(a,h)anthracene	1.000	1.043	-4.3	88	-0.03
37		Benzo(g,h,i)perylene	1.000	1.040	-4.0	87	-0.03

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