Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE091715\
Data File : BE090857.D
Acq On : 17 Sep 2015 9:41
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 17 11:34:18 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 02 11:40:50 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	73	0.00
2		1,4-Dioxane	1.000	0.945	5.5	65	-0.01
3		n-Nitrosodimethylamine	1.000	0.954	4.6	67	0.00
4	S	2-Fluorophenol	1.000	1.072	-7.2	77	0.00
5	S	Phenol-d6	1.000	1.010	-1.0	77	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	82	0.00
7	S	Nitrobenzene-d5	1.000	0.916	8.4	75	0.00
8		Nitrobenzene	1.000	0.902	9.8	77	0.00
9		Naphthalene	1.000	0.947	5.3	78	-0.02
10		Hexachlorobutadiene	1.000	0.908	9.2	72	0.00
11		2-Methylnaphthalene	1.000	1.038	-3.8	88	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	93	0.00
13	S	2,4,6-Tribromophenol	1.000	0.977	2.3	95	0.00
14	S	2-Fluorobiphenyl	1.000	0.901	9.9	85	0.00
15		Acenaphthylene	1.000	0.953	4.7	92	0.00
16		Acenaphthene	1.000	0.937	6.3	86	0.00
17		Fluorene	1.000	0.960	4.0	90	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	86	0.00
19		4-Bromophenyl-phenylether	1.000	0.958	4.2	86	-0.02
20		Hexachlorobenzene	1.000	0.971	2.9	82	0.00
21		Pentachlorophenol	1.000	1.014	-1.4	74	0.00
22		Phenanthrene	1.000	0.973	2.7	84	-0.02
23		Anthracene	1.000	0.993	0.7	89	0.00
24		Fluoranthene	1.000	0.942	5.8	88	-0.01
25	I	Chrysene-d12	5.000	5.000	0.0	82	0.00
26		Pyrene	1.000	1.035	-3.5	89	-0.01
27	S	Terphenyl-d14	1.000	1.019	-1.9	87	0.00
28		Benzo(a)anthracene	1.000	0.980	2.0	81	0.00
29		Chrysene	1.000	1.001	-0.1	79	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.962	3.8	88	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.917	8.3	81	-0.02
32	I	Perylene-d12	5.000	5.000	0.0	82	-0.01
33		Benzo(b)fluoranthene	1.000	0.979	2.1	83	0.00
34		Benzo(k)fluoranthene	1.000	0.947	5.3	79	0.00
35	C	Benzo(a)pyrene	1.000	0.972	2.8	85	-0.01
36		Dibenzo(a,h)anthracene	1.000	0.921	7.9	80	-0.02
37		Benzo(g,h,i)perylene	1.000	0.946	5.4	82	-0.03

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