Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE090115\
Data File : BE090688.D
Acq On : 2 Sep 2015 4:25
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 02 12:25:22 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	90	0.00
2		1,4-Dioxane	1.000	1.117	-11.7	93	0.00
3		n-Nitrosodimethylamine	1.000	0.998	0.2	87	0.00
4	S	2-Fluorophenol	1.000	0.996	0.4	88	0.00
5	S	Phenol-d6	1.000	0.947	5.3	90	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	92	0.00
7	S	Nitrobenzene-d5	1.000	0.964	3.6	89	0.00
8		Nitrobenzene	1.000	0.926	7.4	89	0.00
9		Naphthalene	1.000	1.001	-0.1	92	0.00
10		Hexachlorobutadiene	1.000	1.018	-1.8	90	0.00
11		2-Methylnaphthalene	1.000	0.984	1.6	93	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	92	0.00
13	S	2,4,6-Tribromophenol	1.000	0.883	11.7	83	0.00
14	S	2-Fluorobiphenyl	1.000	0.987	1.3	92	0.00
15		Acenaphthylene	1.000	0.982	1.8	94	0.00
16		Acenaphthene	1.000	1.020	-2.0	93	0.00
17		Fluorene	1.000	0.977	2.3	91	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	89	0.00
19		4-Bromophenyl-phenylether	1.000	0.987	1.3	91	0.00
20		Hexachlorobenzene	1.000	1.062	-6.2	92	0.00
21		Pentachlorophenol	1.000	1.016	-1.6	77	0.00
22		Phenanthrene	1.000	1.013	-1.3	90	0.00
23		Anthracene	1.000	0.976	2.4	90	0.00
24		Fluoranthene	1.000	0.968	3.2	93	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	96	0.00
26		Pyrene	1.000	0.956	4.4	97	0.00
27	S	Terphenyl-d14	1.000	0.940	6.0	94	0.00
28		Benzo(a)anthracene	1.000	1.010	-1.0	98	0.00
29		Chrysene	1.000	1.049	-4.9	97	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.799	20.1	84	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.926	7.4	96	0.00
32	I	Perylene-d12	5.000	5.000	0.0	96	0.00
33		Benzo(b)fluoranthene	1.000	0.972	2.8	96	0.00
34		Benzo(k)fluoranthene	1.000	1.020	-2.0	100	0.00
35	C	Benzo(a)pyrene	1.000	0.963	3.7	99	0.00
36		Dibenzo(a,h)anthracene	1.000	0.941	5.9	96	0.00
37		Benzo(g,h,i)perylene	1.000	0.943	5.7	96	0.00

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