Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE071715\
Data File : BE090167.D
Acq On : 18 Jul 2015 4:46
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 18 07:01:31 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE071715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Jul 18 00:40:53 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	80	0.00
2		1,4-Dioxane	1.000	1.175	-17.5	87	0.00
3		n-Nitrosodimethylamine	1.000	1.034	-3.4	85	0.00
4	S	2-Fluorophenol	1.000	0.899	10.1	73	0.00
5	S	Phenol-d6	1.000	0.925	7.5	74	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	77	0.00
7	S	Nitrobenzene-d5	1.000	0.945	5.5	75	0.00
8		Nitrobenzene	1.000	0.972	2.8	74	0.00
9		Naphthalene	1.000	0.980	2.0	78	0.00
10		Hexachlorobutadiene	1.000	0.943	5.7	74	0.00
11		2-Methylnaphthalene	1.000	0.934	6.6	74	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	74	0.00
13	S	2,4,6-Tribromophenol	1.000	0.730	27.0#	56	0.00
14	S	2-Fluorobiphenyl	1.000	0.974	2.6	73	0.00
15		Acenaphthylene	1.000	0.963	3.7	73	0.00
16		Acenaphthene	1.000	0.957	4.3	73	0.00
17		Fluorene	1.000	0.948	5.2	72	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	72	0.00
19		4-Bromophenyl-phenylether	1.000	0.957	4.3	71	0.00
20		Hexachlorobenzene	1.000	0.963	3.7	71	0.00
21		Pentachlorophenol	1.000	0.787	21.3	56	0.00
22		Phenanthrene	1.000	1.004	-0.4	74	0.00
23		Anthracene	1.000	0.973	2.7	72	0.02
24		Fluoranthene	1.000	0.964	3.6	71	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	72	0.00
26		Pyrene	1.000	0.958	4.2	71	0.00
27	S	Terphenyl-d14	1.000	0.949	5.1	70	0.00
28		Benzo(a)anthracene	1.000	0.917	8.3	69	0.00
29		Chrysene	1.000	0.969	3.1	72	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.658	34.2#	49	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.863	13.7	64	0.00
32	I	Perylene-d12	5.000	5.000	0.0	70	0.00
33		Benzo(b)fluoranthene	1.000	0.930	7.0	67	0.00
34		Benzo(k)fluoranthene	1.000	0.938	6.2	68	0.00
35	C	Benzo(a)pyrene	1.000	0.891	10.9	64	0.00
36		Dibenzo(a,h)anthracene	1.000	0.866	13.4	63	0.00
37		Benzo(g,h,i)perylene	1.000	0.906	9.4	65	0.00

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