Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE092115\
Data File : BE090917.D
Acq On : 21 Sep 2015 14:21
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 22 02:16:22 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE091815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 18 18:17:34 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	128	0.00
2		1,4-Dioxane	1.000	0.961	3.9	113	0.00
3		n-Nitrosodimethylamine	1.000	0.811	18.9	99	0.00
4	S	2-Fluorophenol	1.000	0.738	26.2#	91	0.00
5	S	Phenol-d6	1.000	0.768	23.2	101	0.01
6		bis(2-Chloroethyl)ether	1.000	1.024	-2.4	133	0.00
7	I	Naphthalene-d8	5.000	5.000	0.0	130	0.00
8	S	Nitrobenzene-d5	1.000	0.856	14.4	112	0.00
9		Nitrobenzene	1.000	0.806	19.4	109	0.00
10		Naphthalene	1.000	0.966	3.4	121	0.00
11		Hexachlorobutadiene	1.000	1.046	-4.6	129	0.00
12		2-Methylnaphthalene	1.000	0.885	11.5	118	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	132	0.00
14	S	2,4,6-Tribromophenol	1.000	0.607	39.3#	83	0.00
15	S	2-Fluorobiphenyl	1.000	0.905	9.5	119	0.00
16		Acenaphthylene	1.000	0.879	12.1	117	0.00
17		Acenaphthene	1.000	0.959	4.1	128	0.00
18		Fluorene	1.000	0.935	6.5	123	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	134	0.00
20		4-Bromophenyl-phenylether	1.000	0.876	12.4	119	0.00
21		Hexachlorobenzene	1.000	1.028	-2.8	130	0.00
22		Pentachlorophenol	1.000	0.509	49.1#	54	0.02
23		Phenanthrene	1.000	0.972	2.8	126	0.00
24		Anthracene	1.000	0.875	12.5	121	0.00
25		Fluoranthene	1.000	0.823	17.7	111	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	117	0.00
27		Pyrene	1.000	0.882	11.8	111	0.00
28	S	Terphenyl-d14	1.000	0.880	12.0	107	0.00
29		Benzo(a)anthracene	1.000	0.854	14.6	105	0.00
30		Chrysene	1.000	0.973	2.7	109	0.00
31		Bis(2-ethylhexyl)phthalate	1.000	0.581	41.9#	63	0.00
32		Indeno(1,2,3-cd)pyrene	1.000	0.748	25.2#	87	0.00
33	I	Perylene-d12	5.000	5.000	0.0	104	0.00
34		Benzo(b)fluoranthene	1.000	0.865	13.5	93	0.00
35		Benzo(k)fluoranthene	1.000	0.914	8.6	96	0.00
36	C	Benzo(a)pyrene	1.000	0.829	17.1	90	0.00
37		Dibenzo(a,h)anthracene	1.000	0.810	19.0	85	0.00
38		Benzo(g,h,i)perylene	1.000	0.825	17.5	87	0.00

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