Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE090415\
Data File : BE090730.D
Acq On : 4 Sep 2015 15:23
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 04 16:04:21 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	82	0.00
2		1,4-Dioxane	1.000	1.024	-2.4	79	0.00
3		n-Nitrosodimethylamine	1.000	0.895	10.5	72	0.00
4	S	2-Fluorophenol	1.000	1.041	-4.1	84	0.00
5	S	Phenol-d6	1.000	1.034	-3.4	90	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	89	0.00
7	S	Nitrobenzene-d5	1.000	0.966	3.4	86	0.00
8		Nitrobenzene	1.000	0.946	5.4	88	0.00
9		Naphthalene	1.000	1.010	-1.0	90	0.00
10		Hexachlorobutadiene	1.000	0.980	2.0	84	0.00
11		2-Methylnaphthalene	1.000	1.128	-12.8	104	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	101	0.00
13	S	2,4,6-Tribromophenol	1.000	1.255	-25.5#	137	0.00
14	S	2-Fluorobiphenyl	1.000	0.996	0.4	102	0.00
15		Acenaphthylene	1.000	1.057	-5.7	111	0.00
16		Acenaphthene	1.000	1.012	-1.2	101	0.00
17		Fluorene	1.000	1.009	-0.9	103	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	101	0.00
19		4-Bromophenyl-phenylether	1.000	1.014	-1.4	107	0.00
20		Hexachlorobenzene	1.000	0.994	0.6	99	0.00
21		Pentachlorophenol	1.000	1.193	-19.3	116	0.00
22		Phenanthrene	1.000	1.017	-1.7	103	-0.02
23		Anthracene	1.000	1.056	-5.6	111	0.00
24		Fluoranthene	1.000	1.060	-6.0	116	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	100	0.00
26		Pyrene	1.000	1.147	-14.7	121	0.00
27	S	Terphenyl-d14	1.000	1.111	-11.1	116	0.00
28		Benzo(a)anthracene	1.000	1.073	-7.3	108	0.00
29		Chrysene	1.000	1.107	-10.7	106	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.285	-28.5#	149	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.142	-14.2	123	0.00
32	I	Perylene-d12	5.000	5.000	0.0	106	0.00
33		Benzo(b)fluoranthene	1.000	1.051	-5.1	114	0.00
34		Benzo(k)fluoranthene	1.000	1.075	-7.5	115	0.00
35	C	Benzo(a)pyrene	1.000	1.105	-10.5	125	0.00
36		Dibenzo(a,h)anthracene	1.000	1.072	-7.2	123	0.00
37		Benzo(g,h,i)perylene	1.000	1.061	-6.1	119	0.00

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