Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE080515\
Data File : BE090376.D
Acq On : 6 Aug 2015 1:18
Operator : TP/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC001

Quant Time: Aug 06 13:07:16 2015
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE080515.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 06 12:28:08 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	116	0.00
2		1,4-Dioxane	1.000	1.242	-24.2	125	0.00
3		n-Nitrosodimethylamine	1.000	1.018	-1.8	119	0.00
4	S	2-Fluorophenol	1.000	1.233	-23.3	154	0.00
5	S	Phenol-d6	1.000	1.055	-5.5	137	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	117	0.00
7	S	Nitrobenzene-d5	1.000	0.994	0.6	125	0.00
8		Nitrobenzene	1.000	0.963	3.7	128	0.00
9		Naphthalene	1.000	1.005	-0.5	118	0.00
10		Hexachlorobutadiene	1.000	1.074	-7.4	115	0.00
11		2-Methylnaphthalene	1.000	1.067	-6.7	130	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	118	0.00
13	S	2,4,6-Tribromophenol	1.000	1.220	-22.0	166	0.00
14	S	2-Fluorobiphenyl	1.000	1.064	-6.4	124	0.00
15		Acenaphthylene	1.000	1.049	-4.9	125	0.00
16		Acenaphthene	1.000	0.998	0.2	118	0.00
17		Fluorene	1.000	1.005	-0.5	117	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	112	0.00
19		4-Bromophenyl-phenylether	1.000	1.078	-7.8	123	0.00
20		Hexachlorobenzene	1.000	1.015	-1.5	113	0.00
21		Pentachlorophenol	1.000	1.335	-33.5#	192	0.00
22		Phenanthrene	1.000	0.995	0.5	112	0.00
23		Anthracene	1.000	1.063	-6.3	122	-0.02
24		Fluoranthene	1.000	1.098	-9.8	127	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	117	0.00
26		Pyrene	1.000	1.078	-7.8	131	0.00
27	S	Terphenyl-d14	1.000	1.075	-7.5	127	0.00
28		Benzo(a)anthracene	1.000	0.998	0.2	130	0.00
29		Chrysene	1.000	1.010	-1.0	115	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.131	-13.1	155	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	0.994	0.6	142	0.00
32	I	Perylene-d12	5.000	5.000	0.0	123	0.00
33		Benzo(b)fluoranthene	1.000	1.080	-8.0	148	0.00
34		Benzo(k)fluoranthene	1.000	0.956	4.4	114	0.00
35	C	Benzo(a)pyrene	1.000	1.047	-4.7	141	0.00
36		Dibenzo(a,h)anthracene	1.000	0.998	0.2	147	0.00
37		Benzo(g,h,i)perylene	1.000	1.063	-6.3	135	0.00

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