Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE080515\
Data File : BE090369.D
Acq On : 5 Aug 2015 21:07
Operator : TP/UM
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE080515

Quant Time: Aug 06 13:01:22 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	102	0.00
2		1,4-Dioxane	1.000	1.225	-22.5	109	0.00
3		n-Nitrosodimethylamine	1.000	0.924	7.6	95	0.00
4	S	2-Fluorophenol	1.000	0.939	6.1	103	0.00
5	S	Phenol-d6	1.000	0.961	3.9	106	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	103	0.00
7	S	Nitrobenzene-d5	1.000	0.978	2.2	108	0.00
8		Nitrobenzene	1.000	0.949	5.1	110	0.00
9		Naphthalene	1.000	1.004	-0.4	104	0.00
10		Hexachlorobutadiene	1.000	1.079	-7.9	101	0.00
11		2-Methylnaphthalene	1.000	1.038	-3.8	111	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	105	0.00
13	S	2,4,6-Tribromophenol	1.000	1.007	-0.7	118	0.00
14	S	2-Fluorobiphenyl	1.000	1.052	-5.2	109	0.00
15		Acenaphthylene	1.000	1.040	-4.0	111	0.00
16		Acenaphthene	1.000	1.002	-0.2	105	0.00
17		Fluorene	1.000	1.044	-4.4	109	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	103	0.00
19		4-Bromophenyl-phenylether	1.000	1.047	-4.7	110	0.00
20		Hexachlorobenzene	1.000	1.038	-3.8	106	0.00
21		Pentachlorophenol	1.000	0.989	1.1	120	0.00
22		Phenanthrene	1.000	0.986	1.4	102	0.00
23		Anthracene	1.000	1.042	-4.2	110	0.00
24		Fluoranthene	1.000	1.039	-3.9	111	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	105	0.00
26		Pyrene	1.000	0.999	0.1	110	0.00
27	S	Terphenyl-d14	1.000	1.009	-0.9	107	0.00
28		Benzo(a)anthracene	1.000	0.948	5.2	111	0.00
29		Chrysene	1.000	1.035	-3.5	107	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.900	10.0	105	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.005	-0.5	130	0.00
32	I	Perylene-d12	5.000	5.000	0.0	107	0.00
33		Benzo(b)fluoranthene	1.000	1.015	-1.5	120	0.00
34		Benzo(k)fluoranthene	1.000	1.016	-1.6	106	0.00
35	C	Benzo(a)pyrene	1.000	1.022	-2.2	119	0.00
36		Dibenzo(a,h)anthracene	1.000	1.022	-2.2	133	0.00
37		Benzo(g,h,i)perylene	1.000	1.070	-7.0	119	0.00

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