Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE080315\
Data File : BE090325.D
Acq On : 4 Aug 2015 00:51
Operator : TP/UM
Sample : SSTDICV001
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE080315

Quant Time: Aug 04 02:31:40 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 04 02:31:31 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	136	0.00
2		1,4-Dioxane	1.000	1.371	-37.1#	150	0.00
3		n-Nitrosodimethylamine	1.000	0.946	5.4	127	0.00
4	S	2-Fluorophenol	1.000	1.012	-1.2	141	0.00
5	S	Phenol-d6	1.000	0.952	4.8	139	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	138	0.00
7	S	Nitrobenzene-d5	1.000	0.970	3.0	139	0.00
8		Nitrobenzene	1.000	0.933	6.7	136	0.00
9		Naphthalene	1.000	0.995	0.5	138	0.00
10		Hexachlorobutadiene	1.000	1.024	-2.4	138	0.00
11		2-Methylnaphthalene	1.000	1.045	-4.5	145	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	142	0.00
13	S	2,4,6-Tribromophenol	1.000	1.019	-1.9	174	0.00
14	S	2-Fluorobiphenyl	1.000	1.025	-2.5	144	0.00
15		Acenaphthylene	1.000	1.014	-1.4	145	0.00
16		Acenaphthene	1.000	0.988	1.2	141	0.00
17		Fluorene	1.000	0.980	2.0	149	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	142	0.00
19		4-Bromophenyl-phenylether	1.000	1.037	-3.7	143	0.00
20		Hexachlorobenzene	1.000	1.029	-2.9	143	0.00
21		Pentachlorophenol	1.000	0.952	4.8	163	0.00
22		Phenanthrene	1.000	1.000	0.0	148	0.00
23		Anthracene	1.000	0.986	1.4	148	0.00
24		Fluoranthene	1.000	1.033	-3.3	153	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	152	0.00
26		Pyrene	1.000	1.040	-4.0	151	0.00
27	S	Terphenyl-d14	1.000	1.051	-5.1	153	0.00
28		Benzo(a)anthracene	1.000	1.032	-3.2	176	0.00
29		Chrysene	1.000	1.218	-21.8	158	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	0.937	6.3	168	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.362	-36.2#	277	0.00
32	I	Perylene-d12	5.000	5.000	0.0	154	0.00
33		Benzo(b)fluoranthene	1.000	1.172	-17.2	226	0.00
34		Benzo(k)fluoranthene	1.000	1.280	-28.0#	184	0.00
35	C	Benzo(a)pyrene	1.000	1.190	-19.0	218	0.00
36		Dibenzo(a,h)anthracene	1.000	1.285	-28.5#	272	0.00
37		Benzo(g,h,i)perylene	1.000	1.216	-21.6	203	0.00

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