Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081015\
Data File : BE090414.D
Acq On : 10 Aug 2015 22:36
Operator : TP/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 11 03:56:53 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	132	0.00
2		1,4-Dioxane	1.000	1.192	-19.2	137	0.00
3		n-Nitrosodimethylamine	1.000	1.089	-8.9	144	0.00
4	S	2-Fluorophenol	1.000	1.566	-56.6#	221	0.00
5	S	Phenol-d6	1.000	1.288	-28.8#	200	0.00
6	I	Naphthalene-d8	5.000	5.000	0.0	138	0.00
7	S	Nitrobenzene-d5	1.000	1.086	-8.6	165	0.00
8		Nitrobenzene	1.000	1.050	-5.0	169	0.00
9		Naphthalene	1.000	1.021	-2.1	142	0.00
10		Hexachlorobutadiene	1.000	1.059	-5.9	134	0.00
11		2-Methylnaphthalene	1.000	1.141	-14.1	165	0.00
12	I	Acenaphthene-d10	5.000	5.000	0.0	148	0.00
13	S	2,4,6-Tribromophenol	1.000	1.769	-76.9#	326	0.00
14	S	2-Fluorobiphenyl	1.000	1.072	-7.2	156	0.00
15		Acenaphthylene	1.000	1.097	-9.7	163	0.00
16		Acenaphthene	1.000	1.008	-0.8	149	0.00
17		Fluorene	1.000	1.033	-3.3	151	0.00
18	I	Phenanthrene-d10	5.000	5.000	0.0	140	0.00
19		4-Bromophenyl-phenylether	1.000	1.063	-6.3	153	0.00
20		Hexachlorobenzene	1.000	0.939	6.1	131	0.00
21		Pentachlorophenol	1.000	1.163	-16.3	202	0.00
22		Phenanthrene	1.000	1.014	-1.4	144	0.00
23		Anthracene	1.000	1.135	-13.5	164	-0.02
24		Fluoranthene	1.000	1.145	-14.5	167	0.00
25	I	Chrysene-d12	5.000	5.000	0.0	141	0.00
26		Pyrene	1.000	1.155	-15.5	170	0.00
27	S	Terphenyl-d14	1.000	1.174	-17.4	168	0.00
28		Benzo(a)anthracene	1.000	1.085	-8.5	171	0.00
29		Chrysene	1.000	1.025	-2.5	142	0.00
30		Bis(2-ethylhexyl)phthalate	1.000	1.611	-61.1#	290	0.00
31		Indeno(1,2,3-cd)pyrene	1.000	1.080	-8.0	192	0.00
32	I	Perylene-d12	5.000	5.000	0.0	152	0.00
33		Benzo(b)fluoranthene	1.000	1.090	-9.0	186	0.00
34		Benzo(k)fluoranthene	1.000	0.965	3.5	143	0.00
35	C	Benzo(a)pyrene	1.000	1.087	-8.7	183	0.00
36		Dibenzo(a,h)anthracene	1.000	1.089	-8.9	205	0.00
37		Benzo(g,h,i)perylene	1.000	1.116	-11.6	176	0.00

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