Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE090415\
Data File : BE090728.D
Acq On : 4 Sep 2015 14:00
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Sep 04 14:33:07 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	97	0.00
2		1,4-Dioxane	0.814	0.727	10.7	89	0.00
3		n-Nitrosodimethylamine	1.033	0.892	13.6	88	0.00
4	S	2-Fluorophenol	1.039	1.156	-11.3	115	0.00
5	S	Phenol-d6	1.392	1.616	-16.1	118	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	106	0.00
7	S	Nitrobenzene-d5	0.330	0.336	-1.8	108	0.00
8		Nitrobenzene	0.365	0.360	1.4	109	0.00
9		Naphthalene	1.091	1.093	-0.2	107	0.00
10		Hexachlorobutadiene	0.171	0.162	5.3	98	0.00
11		2-Methylnaphthalene	0.590	0.677	-14.7	126	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	118	0.00
13	S	2,4,6-Tribromophenol	0.143	0.173	-21.0	170#	0.00
14	S	2-Fluorobiphenyl	1.385	1.397	-0.9	121	0.00
15		Acenaphthylene	7.992	8.582	-7.4	132	0.00
16		Acenaphthene	1.335	1.353	-1.3	119	0.00
17		Fluorene	1.666	1.676	-0.6	120	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	111	0.00
19		4-Bromophenyl-phenylether	0.170	0.189	-11.2	128	0.00
20		Hexachlorobenzene	0.928	0.917	1.2	111	0.00
21		Pentachlorophenol	0.054	0.068	-25.9#	159#	0.00
22		Phenanthrene	1.237	1.305	-5.5	117	-0.02
23		Anthracene	1.035	1.159	-12.0	129	0.00
24		Fluoranthene	1.304	1.428	-9.5	131	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	107	0.00
26		Pyrene	1.035	1.242	-20.0	136	0.00
27	S	Terphenyl-d14	0.553	0.641	-15.9	130	0.00
28		Benzo(a)anthracene	1.156	1.213	-4.9	119	0.00
29		Chrysene	1.279	1.342	-4.9	114	0.00
30		Bis(2-ethylhexyl)phthalate	0.345	0.451	-30.7#	191#	0.00
31		Indeno(1,2,3-cd)pyrene	1.181	1.351	-14.4	133	0.00
32	I	Perylene-d12	1.000	1.000	0.0	116	0.00
33		Benzo(b)fluoranthene	1.084	1.134	-4.6	124	0.00
34		Benzo(k)fluoranthene	1.263	1.375	-8.9	128	0.00
35	C	Benzo(a)pyrene	0.970	1.112	-14.6	142	0.00
36		Dibenzo(a,h)anthracene	1.020	1.124	-10.2	133	0.00
37		Benzo(g,h,i)perylene	1.116	1.186	-6.3	130	-0.01

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