Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081415\
Data File : BE090514.D
Acq On : 14 Aug 2015 10:06
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Aug 14 15:26:29 2015
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE081415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 14 15:17:14 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	116	0.00
2		1,4-Dioxane	0.658	0.685	-4.1	115	0.00
3		n-Nitrosodimethylamine	0.863	0.908	-5.2	124	0.00
4	S	2-Fluorophenol	1.446	1.468	-1.5	118	0.00
5	S	Phenol-d6	2.125	2.152	-1.3	123	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	117	0.00
7	S	Nitrobenzene-d5	0.398	0.400	-0.5	123	0.00
8		Nitrobenzene	0.413	0.414	-0.2	125	0.00
9		Naphthalene	1.136	1.134	0.2	117	0.00
10		Hexachlorobutadiene	0.167	0.166	0.6	116	0.00
11		2-Methylnaphthalene	0.740	0.759	-2.6	123	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	121	0.00
13	S	2,4,6-Tribromophenol	0.192	0.191	0.5	133	-0.02
14	S	2-Fluorobiphenyl	1.493	1.501	-0.5	122	0.00
15		Acenaphthylene	9.132	9.178	-0.5	124	0.00
16		Acenaphthene	1.374	1.366	0.6	122	0.00
17		Fluorene	1.684	1.704	-1.2	123	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	122	0.00
19		4-Bromophenyl-phenylether	0.204	0.208	-2.0	127	0.00
20		Hexachlorobenzene	0.906	0.890	1.8	120	0.00
21		Pentachlorophenol	0.079	0.074	6.3	152#	0.00
22		Phenanthrene	1.312	1.282	2.3	121	0.00
23		Anthracene	1.208	1.213	-0.4	127	0.00
24		Fluoranthene	1.396	1.425	-2.1	128	0.00
25	I	Chrysene-d12	1.000	1.000	0.0	126	0.00
26		Pyrene	1.430	1.409	1.5	128	0.00
27	S	Terphenyl-d14	0.914	0.912	0.2	128	0.00
28		Benzo(a)anthracene	1.316	1.281	2.7	129	0.00
29		Chrysene	1.344	1.352	-0.6	128	0.00
30		Bis(2-ethylhexyl)phthalate	0.725	0.750	-3.4	159#	0.00
31		Indeno(1,2,3-cd)pyrene	1.266	1.245	1.7	135	0.00
32	I	Perylene-d12	1.000	1.000	0.0	129	0.00
33		Benzo(b)fluoranthene	1.184	1.210	-2.2	135	0.00
34		Benzo(k)fluoranthene	1.342	1.355	-1.0	129	0.00
35	C	Benzo(a)pyrene	1.108	1.114	-0.5	134	0.00
36		Dibenzo(a,h)anthracene	1.069	1.057	1.1	135	0.00
37		Benzo(g,h,i)perylene	1.169	1.169	0.0	131	0.00

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