Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE100815\
Data File : BE090996.D
Acq On : 8 Oct 2015 10:36
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Oct 09 03:05:01 2015
Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 07 17:11:43 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	172	0.00
2		1,4-Dioxane	1.000	2.622	-162.2#	398	-0.07
3		n-Nitrosodimethylamine	1.000	0.689	31.1#	123	-0.10
4	S	2-Fluorophenol	1.000	1.147	-14.7	207	-0.02
5	S	Phenol-d6	1.000	1.062	-6.2	204	-0.02
6		bis(2-Chloroethyl)ether	1.000	2.676	-167.6#	424	-0.04
7	I	Naphthalene-d8	5.000	5.000	0.0	173	0.00
8	S	Nitrobenzene-d5	1.000	1.264	-26.4#	251	-0.01
9		Nitrobenzene	1.000	1.324	-32.4#	277	0.00
10		Naphthalene	1.000	0.996	0.4	172	0.00
11		Hexachlorobutadiene	1.000	0.826	17.4	142	0.00
12		2-Methylnaphthalene	1.000	1.078	-7.8	198	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	173	0.00
14	S	2,4,6-Tribromophenol	1.000	1.014	-1.4	218	0.00
15	S	2-Fluorobiphenyl	1.000	0.987	1.3	184	0.00
16		Acenaphthylene	1.000	0.983	1.7	186	0.00
17		Acenaphthene	1.000	0.945	5.5	174	0.00
18		Fluorene	1.000	0.981	1.9	182	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	177	-0.01
20		4-Bromophenyl-phenylether	1.000	0.985	1.5	189	-0.01
21		Hexachlorobenzene	1.000	0.125	87.5#	39	0.00
22		Pentachlorophenol	1.000	1.180	-18.0	258	0.00
23		Phenanthrene	1.000	0.982	1.8	181	-0.01
24		Anthracene	1.000	1.025	-2.5	195	-0.01
25		Fluoranthene	1.000	1.023	-2.3	197	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	186	0.00
27		Benzidine	1.000	2.139	-113.9#	487	-0.02
28		Pyrene	1.000	0.961	3.9	195	0.00
29	S	Terphenyl-d14	1.000	1.034	-3.4	211	0.00
30		Benzo(a)anthracene	1.000	1.066	-6.6	208	0.00
31		3,3'-Dichlorobenzidine	1.000	0.022	97.8#	0	-20.84#
32		Chrysene	1.000	0.913	8.7	177	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	1.405	-40.5#	335	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	1.160	-16.0	239	0.00
35	I	Perylene-d12	5.000	5.000	0.0	223	0.00
36		Benzo(b)fluoranthene	1.000	1.118	-11.8	278	0.00
37		Benzo(k)fluoranthene	1.000	0.785	21.5	188	0.00
38	C	Benzo(a)pyrene	1.000	1.084	-8.4	280	0.00
39		Dibenzo(a,h)anthracene	1.000	1.241	-24.1	329	0.00
40		Benzo(g,h,i)perylene	1.000	1.174	-17.4	302	0.00

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Compound	Amount	Calc.	%Dev	Area%	Dev(min)

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