Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA E\DATA\BE100815\
Data File : BE090998.D
Acq On : 8 Oct 2015 11:57
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Oct 09 03:17:49 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	167	0.00
2		1,4-Dioxane	1.000	2.740	-174.0#	402	-0.10
3		n-Nitrosodimethylamine	1.000	0.832	16.8	143	-0.13
4	S	2-Fluorophenol	1.000	1.182	-18.2	207	-0.02
5	S	Phenol-d6	1.000	1.149	-14.9	215	-0.02
6		bis(2-Chloroethyl)ether	1.000	2.765	-176.5#	425	-0.05
7	I	Naphthalene-d8	5.000	5.000	0.0	172	0.00
8	S	Nitrobenzene-d5	1.000	1.262	-26.2#	249	-0.02
9		Nitrobenzene	1.000	1.363	-36.3#	286	-0.01
10		Naphthalene	1.000	1.002	-0.2	172	-0.02
11		Hexachlorobutadiene	1.000	0.837	16.3	143	-0.02
12		2-Methylnaphthalene	1.000	1.084	-8.4	198	0.00
13	I	Acenaphthene-d10	5.000	5.000	0.0	176	0.00
14	S	2,4,6-Tribromophenol	1.000	1.083	-8.3	240	0.00
15	S	2-Fluorobiphenyl	1.000	0.976	2.4	185	-0.01
16		Acenaphthylene	1.000	0.989	1.1	190	0.00
17		Acenaphthene	1.000	0.949	5.1	178	0.00
18		Fluorene	1.000	0.984	1.6	186	0.00
19	I	Phenanthrene-d10	5.000	5.000	0.0	185	-0.01
20		4-Bromophenyl-phenylether	1.000	0.963	3.7	194	-0.01
21		Hexachlorobenzene	1.000	0.118	88.2#	40	-0.01
22		Pentachlorophenol	1.000	1.307	-30.7#	302	0.00
23		Phenanthrene	1.000	0.961	3.9	186	-0.01
24		Anthracene	1.000	1.020	-2.0	204	-0.01
25		Fluoranthene	1.000	1.029	-2.9	208	0.00
26	I	Chrysene-d12	5.000	5.000	0.0	195	0.00
27		Benzidine	1.000	2.263	-126.3#	552	-0.03
28		Pyrene	1.000	0.978	2.2	207	0.00
29	S	Terphenyl-d14	1.000	1.038	-3.8	221	0.00
30		Benzo(a)anthracene	1.000	1.081	-8.1	221	0.00
31		3,3'-Dichlorobenzidine	1.000	0.022	97.8#	0	-20.84#
32		Chrysene	1.000	0.931	6.9	188	0.00
33		Bis(2-ethylhexyl)phthalate	1.000	1.491	-49.1#	375	0.00
34		Indeno(1,2,3-cd)pyrene	1.000	1.119	-11.9	240	0.00
35	I	Perylene-d12	5.000	5.000	0.0	229	0.00
36		Benzo(b)fluoranthene	1.000	1.126	-12.6	288	0.00
37		Benzo(k)fluoranthene	1.000	0.792	20.8	195	0.00
38	C	Benzo(a)pyrene	1.000	1.088	-8.8	289	0.00
39		Dibenzo(a,h)anthracene	1.000	1.201	-20.1	327	0.00
40		Benzo(g,h,i)perylene	1.000	1.139	-13.9	301	0.00

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

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