Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE102315\
Data File : BE091045.D
Acq On : 23 Oct 2015 22:08
Operator : UM/NP
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Oct 24 03:33:16 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE092815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Oct 24 03:31:05 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	47	-0.05
2		1,4-Dioxane	1.000	0.872	12.8	42	-0.18
3		n-Nitrosodimethylamine	1.000	0.963	3.7	46	-0.27
4	S	2-Fluorophenol	1.000	1.367	-36.7#	67	-0.07
5	S	Phenol-d6	1.000	1.364	-36.4#	71	-0.07
6		bis(2-Chloroethyl)ether	1.000	1.806	-80.6#	77	-0.11
7	I	Naphthalene-d8	5.000	5.000	0.0	48	-0.05
8	S	Nitrobenzene-d5	1.000	1.413	-41.3#	79	-0.06
9		Nitrobenzene	1.000	1.400	-40.0#	82	-0.06
10		Naphthalene	1.000	1.012	-1.2	48	-0.06
11		Hexachlorobutadiene	1.000	0.842	15.8	40	-0.05
12		2-Methylnaphthalene	1.000	1.095	-9.5	55	-0.06
13	I	Acenaphthene-d10	5.000	5.000	0.0	51	-0.05
14	S	2,4,6-Tribromophenol	1.000	1.314	-31.4#	87	-0.05
15	S	2-Fluorobiphenyl	1.000	0.944	5.6	52	-0.06
16		Acenaphthylene	1.000	1.021	-2.1	57	-0.05
17		Acenaphthene	1.000	0.965	3.5	53	-0.05
18		Fluorene	1.000	1.022	-2.2	56	-0.05
19	I	Phenanthrene-d10	5.000	5.000	0.0	57	-0.05
20		4-Bromophenyl-phenylether	1.000	0.917	8.3	57	-0.05
21		Hexachlorobenzene	1.000	0.106	89.4#	12	-0.05
22		Pentachlorophenol	1.000	1.474	-47.4#	106	-0.04
23		Phenanthrene	1.000	0.998	0.2	59	-0.05
24		Anthracene	1.000	1.057	-5.7	65	-0.05
25		Fluoranthene	1.000	1.089	-8.9	68	-0.05
26	I	Chrysene-d12	5.000	5.000	0.0	60	-0.05
27		Benzidine	1.000	2.260	-126.0#	170	-0.08
28		Pyrene	1.000	1.031	-3.1	67	-0.05
29	S	Terphenyl-d14	1.000	1.201	-20.1	79	-0.05
30		Benzo(a)anthracene	1.000	1.091	-9.1	69	-0.05
31		3,3'-Dichlorobenzidine	1.000	0.022	97.8#	0	-20.84#
32		Chrysene	1.000	0.923	7.7	58	-0.04
33		Bis(2-ethylhexyl)phthalate	1.000	1.497	-49.7#	116	-0.05
34		Indeno(1,2,3-cd)pyrene	1.000	1.055	-5.5	69	-0.08
35	I	Perylene-d12	5.000	5.000	0.0	66	-0.06
36		Benzo(b)fluoranthene	1.000	1.109	-10.9	82	-0.05
37		Benzo(k)fluoranthene	1.000	0.839	16.1	59	-0.05
38	C	Benzo(a)pyrene	1.000	1.102	-10.2	85	-0.05
39		Dibenzo(a,h)anthracene	1.000	1.179	-17.9	93	-0.08
40		Benzo(g,h,i)perylene	1.000	1.134	-13.4	87	-0.09

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

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