Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE062217\
Data File : BE093325.D
Acq On : 22 Jun 2017 12:23
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 23 07:50:06 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE061217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 12 14:07:54 2017
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	0.400	0.400	0.0	31	-0.02
2		1,4-Dioxane	0.400	0.367	8.3	34	0.00
3		n-Nitrosodimethylamine	0.400	0.422	-5.5	36	0.00
4	S	2-Fluorophenol	0.400	0.502	-25.5#	45	0.00
5	S	Phenol-d6	0.400	0.513	-28.2#	45	0.00
6		bis(2-Chloroethyl)ether	0.400	0.446	-11.5	39	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	36	-0.02
8	S	Nitrobenzene-d5	0.400	0.485	-21.2	50	0.00
9		Naphthalene	0.400	0.430	-7.5	44	0.00
10		Hexachlorobutadiene	0.400	0.413	-3.2	43	-0.02
11	SURR	2-Methylnaphthalene-d10	0.400	0.466	-16.5	49	-0.01
12		2-Methylnaphthalene	0.400	0.450	-12.5	46	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	41	0.00
14	S	2,4,6-Tribromophenol	0.400	0.504	-26.0#	63	0.00
15	S	2-Fluorobiphenyl	0.400	0.409	-2.2	47	0.00
16		Acenaphthylene	0.400	0.447	-11.7	55	0.00
17		Acenaphthene	0.400	0.418	-4.5	50	-0.02
18		Fluorene	0.400	0.421	-5.2	50	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	33	-0.03
20		4-Bromophenyl-phenylether	0.400	0.316	21.0	30	0.00
21		Hexachlorobenzene	0.400	0.507	-26.7#	54	0.00
22		Pentachlorophenol	0.400	0.474	-18.5	49	0.00
23		Phenanthrene	0.400	0.384	4.0	37	0.00
24		Anthracene	0.400	0.452	-13.0	48	0.00
25	SURR	Fluoranthene-d10	0.400	0.414	-3.5	40	0.00
26		Fluoranthene	0.400	0.407	-1.7	41	-0.02
27	I	Chrysene-d12	0.400	0.400	0.0	28	-0.02
28		Pyrene	0.400	0.454	-13.5	35	0.00
29	S	Terphenyl-d14	0.400	0.466	-16.5	36	0.00
30		Benzo(a)anthracene	0.400	0.448	-12.0	36	0.00
31		Chrysene	0.400	0.419	-4.7	33	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.493	-23.2	37	-0.02
33		Indeno(1,2,3-cd)pyrene	0.400	0.425	-6.2	34	0.00
34	I	Perylene-d12	0.400	0.400	0.0	27	0.00
35		Benzo(b)fluoranthene	0.400	0.467	-16.8	37	0.00
36		Benzo(k)fluoranthene	0.400	0.406	-1.5	31	0.00
37	C	Benzo(a)pyrene	0.400	0.448	-12.0	35	0.00
38		Dibenzo(a,h)anthracene	0.400	0.426	-6.5	33	0.00
39		Benzo(g,h,i)perylene	0.400	0.421	-5.2	33	0.02

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

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