Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE110217\
Data File : BE094699.D
Acq On : 2 Nov 2017 17:27
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Nov 02 18:21:11 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE102317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Oct 23 22:40:47 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	115	0.00
2		1,4-Dioxane	0.400	0.384	4.0	95	-0.01
3		n-Nitrosodimethylamine	0.400	0.292	27.0#	84	0.01
4	S	2-Fluorophenol	0.400	0.297	25.8#	85	0.05
5	S	Phenol-d6	0.400	0.359	10.3	106	0.04
6		bis(2-Chloroethyl)ether	0.400	0.353	11.8	102	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	121	0.00
8	S	Nitrobenzene-d5	0.400	0.337	15.8	106	0.03
9		Naphthalene	0.400	0.352	12.0	103	0.02
10		Hexachlorobutadiene	0.400	0.382	4.5	114	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.346	13.5	102	0.00
12		2-Methylnaphthalene	0.400	0.343	14.2	103	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	108	0.01
14	S	2,4,6-Tribromophenol	0.400	0.262	34.5#	74	0.03
15	S	2-Fluorobiphenyl	0.400	0.364	9.0	97	0.01
16		Acenaphthylene	0.400	0.344	14.0	94	0.00
17		Acenaphthene	0.400	0.358	10.5	96	0.00
18		Fluorene	0.400	0.342	14.5	93	0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	102	0.03
20		4-Bromophenyl-phenylether	0.400	0.329	17.8	86	0.03
21		Hexachlorobenzene	0.400	0.571	-42.7#	127	0.00
22		Pentachlorophenol	0.400	0.324	19.0	83	0.07
23		Phenanthrene	0.400	0.511	-27.7#	118	0.00
24		Anthracene	0.400	0.395	1.3	99	0.00
25	SURR	Fluoranthene-d10	0.400	0.411	-2.7	96	0.00
26		Fluoranthene	0.400	0.402	-0.5	95	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	105	0.00
28		Pyrene	0.400	0.356	11.0	95	0.00
29	S	Terphenyl-d14	0.400	0.349	12.8	93	0.00
30		Benzo(a)anthracene	0.400	0.350	12.5	92	0.01
31		Chrysene	0.400	0.365	8.8	99	0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.331	17.3	95	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.271	32.3#	72	0.05
34	I	Perylene-d12	0.400	0.400	0.0	99	0.01
35		Benzo(b)fluoranthene	0.400	0.359	10.3	91	0.00
36		Benzo(k)fluoranthene	0.400	0.398	0.5	99	0.01
37	C	Benzo(a)pyrene	0.400	0.373	6.8	93	0.02
38		Dibenzo(a,h)anthracene	0.400	0.305	23.8	77	0.03
39		Benzo(g,h,i)perylene	0.400	0.266	33.5#	67	0.06

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

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