Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE121119\
Data File : BE100857.D
Acq On : 11 Dec 2019 23:03
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 12 04:43:44 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE120419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 04 17:18:11 2019
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	93	-0.01
2		1,4-Dioxane	0.400	0.446	-11.5	104	-0.01
3		n-Nitrosodimethylamine	0.400	0.412	-3.0	110	-0.01
4	S	2-Fluorophenol	0.400	0.500	-25.0	124	-0.01
5	S	Phenol-d6	0.400	0.529	-32.3#	133	0.00
6		bis(2-Chloroethyl)ether	0.400	0.461	-15.3	113	-0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	104	0.00
8	S	Nitrobenzene-d5	0.400	0.554	-38.5#	166	0.00
9		Naphthalene	0.400	0.453	-13.2	123	0.00
10		Hexachlorobutadiene	0.400	0.440	-10.0	118	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.455	-13.7	126	0.00
12		2-Methylnaphthalene	0.400	0.468	-17.0	129	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	113	0.00
14	S	2,4,6-Tribromophenol	0.400	0.654	-63.5#	211	-0.01
15	S	2-Fluorobiphenyl	0.400	0.413	-3.2	122	-0.01
16		Acenaphthylene	0.400	0.441	-10.2	144	0.00
17		Acenaphthene	0.400	0.449	-12.2	139	0.00
18		Fluorene	0.400	0.440	-10.0	139	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	105	-0.01
20		4-Bromophenyl-phenylether	0.400	0.443	-10.7	128	-0.01
21		Hexachlorobenzene	0.400	0.465	-16.3	130	0.00
22		Pentachlorophenol	0.400	0.878	-119.5#	1189	-0.01
23		Phenanthrene	0.400	0.457	-14.2	127	0.00
24		Anthracene	0.400	0.420	-5.0	128	0.00
25	SURR	Fluoranthene-d10	0.400	0.373	6.8	108	0.00
26		Fluoranthene	0.400	0.389	2.8	113	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	84	0.00
28		Pyrene	0.400	0.503	-25.7#	113	0.00
29	S	Terphenyl-d14	0.400	0.454	-13.5	100	0.00
30		Benzo(a)anthracene	0.400	0.457	-14.2	108	-0.01
31		Chrysene	0.400	0.484	-21.0	105	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.598	-49.5#	130	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.495	-23.7	118	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	97	-0.01
35		Benzo(b)fluoranthene	0.400	0.429	-7.2	110	-0.01
36		Benzo(k)fluoranthene	0.400	0.430	-7.5	110	-0.01
37	C	Benzo(a)pyrene	0.400	0.413	-3.2	110	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.433	-8.2	119	-0.03
39		Benzo(g,h,i)perylene	0.400	0.435	-8.7	112	-0.02

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

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