Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN092719\
Data File : BN007888.D
Acq On : 27 Sep 2019 16:40
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Sep 27 18:45:57 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN091819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 18 18:24:40 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	166	-0.02
2		1,4-Dioxane	0.400	0.455	-13.7	223	-0.02
3		n-Nitrosodimethylamine	0.400	0.707	-76.7#	193	-0.02
4	S	2-Fluorophenol	0.400	0.380	5.0	181	-0.02
5	S	Phenol-d6	0.400	0.407	-1.7	186	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.517	-29.3#	243	-0.03
7	I	Naphthalene-d8	0.400	0.400	0.0	168	-0.01
8	S	Nitrobenzene-d5	0.400	0.446	-11.5	193	-0.01
9		Naphthalene	0.400	0.427	-6.7	180	-0.03
10		Hexachlorobutadiene	0.400	0.440	-10.0	184	-0.03
11	SURR	2-Methylnaphthalene-d10	0.400	0.424	-6.0	184	-0.02
12		2-Methylnaphthalene	0.400	0.430	-7.5	186	-0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	161	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.388	3.0	186	-0.01
15	S	2-Fluorobiphenyl	0.400	0.406	-1.5	176	-0.01
16		Acenaphthylene	0.400	0.428	-7.0	191	-0.02
17		Acenaphthene	0.400	0.397	0.8	174	-0.02
18		Fluorene	0.400	0.405	-1.3	177	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	146	-0.01
20		4-Bromophenyl-phenylether	0.400	0.437	-9.2	162	-0.02
21		Hexachlorobenzene	0.400	0.442	-10.5	163	-0.01
22		Pentachlorophenol	0.400	0.534	-33.5#	224	-0.02
23		Phenanthrene	0.400	0.432	-8.0	160	-0.01
24		Anthracene	0.400	0.451	-12.7	171	-0.01
25	SURR	Fluoranthene-d10	0.400	0.417	-4.2	154	-0.01
26		Fluoranthene	0.400	0.421	-5.2	156	-0.01
27	I	Chrysene-d12	0.400	0.400	0.0	137	-0.01
28		Pyrene	0.400	0.451	-12.7	160	-0.01
29	S	Terphenyl-d14	0.400	0.448	-12.0	159	-0.01
30		Benzo(a)anthracene	0.400	0.430	-7.5	153	-0.01
31		Chrysene	0.400	0.422	-5.5	148	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.733	-83.2#	280	-0.02
33		Indeno(1,2,3-cd)pyrene	0.400	0.391	2.3	137	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	133	-0.01
35		Benzo(b)fluoranthene	0.400	0.429	-7.2	151	-0.01
36		Benzo(k)fluoranthene	0.400	0.412	-3.0	142	-0.01
37	C	Benzo(a)pyrene	0.400	0.421	-5.2	147	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.391	2.3	136	-0.02
39		Benzo(g,h,i)perylene	0.400	0.406	-1.5	140	-0.02

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

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