Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121719\
Data File : BN009018.D
Acq On : 17 Dec 2019 18:54
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Dec 18 12:41:09 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN120919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 09 17:02:53 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	137	-0.02
2		1,4-Dioxane	0.400	0.380	5.0	110	0.00
3		n-Nitrosodimethylamine	0.400	0.387	3.3	137	0.00
4	S	2-Fluorophenol	0.400	0.621	-55.2#	197	0.00
5	S	Phenol-d6	0.400	0.531	-32.8#	171	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.418	-4.5	129	-0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	151	-0.01
8	S	Nitrobenzene-d5	0.400	0.400	0.0	146	-0.01
9		Naphthalene	0.400	0.423	-5.7	143	-0.01
10		Hexachlorobutadiene	0.400	0.424	-6.0	143	-0.03
11	SURR	2-Methylnaphthalene-d10	0.400	0.431	-7.7	146	-0.02
12		2-Methylnaphthalene	0.400	0.439	-9.7	147	-0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	153	0.00
14	S	2,4,6-Tribromophenol	0.400	0.379	5.3	139	-0.02
15	S	2-Fluorobiphenyl	0.400	0.398	0.5	143	-0.02
16		Acenaphthylene	0.400	0.447	-11.7	154	-0.02
17		Acenaphthene	0.400	0.427	-6.7	145	-0.02
18		Fluorene	0.400	0.419	-4.7	141	-0.02
19	I	Phenanthrene-d10	0.400	0.400	0.0	151	-0.01
20		4-Bromophenyl-phenylether	0.400	0.404	-1.0	140	-0.01
21		Hexachlorobenzene	0.400	0.413	-3.2	142	-0.01
22		Pentachlorophenol	0.400	0.254	36.5#	99	0.00
23		Phenanthrene	0.400	0.404	-1.0	139	-0.01
24		Anthracene	0.400	0.415	-3.7	145	0.00
25	SURR	Fluoranthene-d10	0.400	0.398	0.5	140	-0.01
26		Fluoranthene	0.400	0.401	-0.3	138	-0.01
27	I	Chrysene-d12	0.400	0.400	0.0	134	0.00
28		Pyrene	0.400	0.485	-21.2	139	0.00
29	S	Terphenyl-d14	0.400	0.457	-14.2	138	0.00
30		Benzo(a)anthracene	0.400	0.442	-10.5	130	0.00
31		Chrysene	0.400	0.423	-5.7	121	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.591	-47.7#	198	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.356	11.0	114	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	122	-0.01
35		Benzo(b)fluoranthene	0.400	0.428	-7.0	111	-0.01
36		Benzo(k)fluoranthene	0.400	0.430	-7.5	115	0.00
37	C	Benzo(a)pyrene	0.400	0.425	-6.2	115	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.408	-2.0	114	-0.02
39		Benzo(g,h,i)perylene	0.400	0.418	-4.5	115	-0.02

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

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