Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121719\
Data File : BN009008.D
Acq On : 17 Dec 2019 10:05
Operator : JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Dec 18 12:21:42 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	119	-0.02
2		1,4-Dioxane	0.400	0.403	-0.8	101	0.00
3		n-Nitrosodimethylamine	0.400	0.356	11.0	109	0.00
4	S	2-Fluorophenol	0.400	0.420	-5.0	115	-0.02
5	S	Phenol-d6	0.400	0.432	-8.0	120	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.422	-5.5	113	-0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	126	-0.01
8	S	Nitrobenzene-d5	0.400	0.396	1.0	120	-0.01
9		Naphthalene	0.400	0.425	-6.2	120	-0.01
10		Hexachlorobutadiene	0.400	0.415	-3.7	117	-0.03
11	SURR	2-Methylnaphthalene-d10	0.400	0.424	-6.0	120	-0.02
12		2-Methylnaphthalene	0.400	0.434	-8.5	121	-0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	124	0.00
14	S	2,4,6-Tribromophenol	0.400	0.401	-0.3	120	-0.02
15	S	2-Fluorobiphenyl	0.400	0.402	-0.5	117	-0.02
16		Acenaphthylene	0.400	0.437	-9.2	122	-0.02
17		Acenaphthene	0.400	0.423	-5.7	117	-0.02
18		Fluorene	0.400	0.421	-5.2	115	-0.02
19	I	Phenanthrene-d10	0.400	0.400	0.0	120	-0.01
20		4-Bromophenyl-phenylether	0.400	0.415	-3.7	114	-0.01
21		Hexachlorobenzene	0.400	0.420	-5.0	115	-0.01
22		Pentachlorophenol	0.400	0.391	2.3	121	-0.01
23		Phenanthrene	0.400	0.414	-3.5	113	-0.01
24		Anthracene	0.400	0.410	-2.5	114	-0.01
25	SURR	Fluoranthene-d10	0.400	0.393	1.8	110	-0.01
26		Fluoranthene	0.400	0.402	-0.5	110	-0.01
27	I	Chrysene-d12	0.400	0.400	0.0	108	0.00
28		Pyrene	0.400	0.477	-19.2	110	0.00
29	S	Terphenyl-d14	0.400	0.449	-12.2	109	0.00
30		Benzo(a)anthracene	0.400	0.437	-9.2	104	0.00
31		Chrysene	0.400	0.423	-5.7	98	-0.02
32		Bis(2-ethylhexyl)phthalate	0.400	0.489	-22.2	132	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.331	17.3	86	-0.02
34	I	Perylene-d12	0.400	0.400	0.0	94	-0.02
35		Benzo(b)fluoranthene	0.400	0.430	-7.5	87	-0.02
36		Benzo(k)fluoranthene	0.400	0.435	-8.7	90	-0.01
37	C	Benzo(a)pyrene	0.400	0.421	-5.2	88	-0.02
38		Dibenzo(a,h)anthracene	0.400	0.393	1.8	85	-0.02
39		Benzo(g,h,i)perylene	0.400	0.404	-1.0	86	-0.02

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

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