Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN070119\
Data File : BN006668.D
Acq On : 01 Jul 2019 19:22
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Jul 02 05:06:50 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN062819.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 28 13:47:56 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	125	0.00
2		1,4-Dioxane	0.400	0.380	5.0	127	0.00
3		n-Nitrosodimethylamine	0.400	0.360	10.0	127	0.00
4	S	2-Fluorophenol	0.400	0.428	-7.0	140	0.00
5	S	Phenol-d6	0.400	0.439	-9.7	145	0.00
6		bis(2-Chloroethyl)ether	0.400	0.376	6.0	113	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	116	-0.01
8	S	Nitrobenzene-d5	0.400	0.431	-7.7	134	-0.01
9		Naphthalene	0.400	0.423	-5.7	121	-0.01
10		Hexachlorobutadiene	0.400	0.430	-7.5	121	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.395	1.3	112	0.00
12		2-Methylnaphthalene	0.400	0.404	-1.0	114	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	97	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.417	-4.2	112	-0.01
15	S	2-Fluorobiphenyl	0.400	0.442	-10.5	107	-0.01
16		Acenaphthylene	0.400	0.458	-14.5	113	0.00
17		Acenaphthene	0.400	0.432	-8.0	102	0.00
18		Fluorene	0.400	0.426	-6.5	101	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	88	-0.01
20		4-Bromophenyl-phenylether	0.400	0.437	-9.2	97	-0.01
21		Hexachlorobenzene	0.400	0.431	-7.7	92	0.00
22		Pentachlorophenol	0.400	0.578	-44.5#	152	-0.02
23		Phenanthrene	0.400	0.421	-5.2	92	0.00
24		Anthracene	0.400	0.425	-6.2	96	0.00
25	SURR	Fluoranthene-d10	0.400	0.408	-2.0	89	0.00
26		Fluoranthene	0.400	0.412	-3.0	88	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	76	-0.02
28		Pyrene	0.400	0.471	-17.7	87	-0.01
29	S	Terphenyl-d14	0.400	0.463	-15.8	88	0.00
30		Benzo(a)anthracene	0.400	0.449	-12.2	87	-0.02
31		Chrysene	0.400	0.431	-7.7	78	-0.02
32		Bis(2-ethylhexyl)phthalate	0.400	0.625	-56.2#	133	-0.02
33		Indeno(1,2,3-cd)pyrene	0.400	0.524	-31.0#	86	-0.05
34	I	Perylene-d12	0.400	0.400	0.0	75	-0.04
35		Benzo(b)fluoranthene	0.400	0.412	-3.0	77	-0.04
36		Benzo(k)fluoranthene	0.400	0.394	1.5	72	-0.04
37	C	Benzo(a)pyrene	0.400	0.429	-7.2	77	-0.04
38		Dibenzo(a,h)anthracene	0.400	0.474	-18.5	84	-0.05
39		Benzo(g,h,i)perylene	0.400	0.493	-23.2	87	-0.05

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

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