Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN070119\
Data File : BN006686.D
Acq On : 02 Jul 2019 07:28
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Jul 02 08:48:16 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	112	0.00
2		1,4-Dioxane	0.400	0.400	0.0	120	0.00
3		n-Nitrosodimethylamine	0.400	0.383	4.3	121	0.00
4	S	2-Fluorophenol	0.400	0.452	-13.0	132	0.00
5	S	Phenol-d6	0.400	0.453	-13.2	134	0.00
6		bis(2-Chloroethyl)ether	0.400	0.383	4.3	103	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	106	-0.01
8	S	Nitrobenzene-d5	0.400	0.437	-9.2	124	-0.01
9		Naphthalene	0.400	0.423	-5.7	110	-0.01
10		Hexachlorobutadiene	0.400	0.430	-7.5	110	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.401	-0.3	103	0.00
12		2-Methylnaphthalene	0.400	0.407	-1.7	104	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	89	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.457	-14.2	113	-0.01
15	S	2-Fluorobiphenyl	0.400	0.443	-10.7	99	-0.01
16		Acenaphthylene	0.400	0.522	-30.5#	118	0.00
17		Acenaphthene	0.400	0.454	-13.5	99	0.00
18		Fluorene	0.400	0.447	-11.7	98	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	82	-0.01
20		4-Bromophenyl-phenylether	0.400	0.443	-10.7	92	-0.01
21		Hexachlorobenzene	0.400	0.427	-6.7	85	0.00
22		Pentachlorophenol	0.400	0.617	-54.2#	157	-0.02
23		Phenanthrene	0.400	0.430	-7.5	87	0.00
24		Anthracene	0.400	0.448	-12.0	95	0.00
25	SURR	Fluoranthene-d10	0.400	0.421	-5.2	86	0.00
26		Fluoranthene	0.400	0.421	-5.2	84	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	72	-0.01
28		Pyrene	0.400	0.480	-20.0	85	0.00
29	S	Terphenyl-d14	0.400	0.469	-17.2	84	0.00
30		Benzo(a)anthracene	0.400	0.467	-16.8	86	-0.02
31		Chrysene	0.400	0.422	-5.5	72	-0.02
32		Bis(2-ethylhexyl)phthalate	0.400	0.703	-75.7#	142	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.523	-30.8#	81	-0.03
34	I	Perylene-d12	0.400	0.400	0.0	72	-0.03
35		Benzo(b)fluoranthene	0.400	0.413	-3.2	74	-0.03
36		Benzo(k)fluoranthene	0.400	0.385	3.8	68	-0.03
37	C	Benzo(a)pyrene	0.400	0.430	-7.5	75	-0.03
38		Dibenzo(a,h)anthracene	0.400	0.475	-18.7	81	-0.03
39		Benzo(g,h,i)perylene	0.400	0.492	-23.0	83	-0.03

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

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