Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\terastorage\svoasrv\HPCHEM1\BNA_N\Data\BN062819\
Data File : BN006618.D
Acq On : 28 Jun 2019 18:08
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Jun 28 18:42:01 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	102	0.00
2		1,4-Dioxane	0.400	0.390	2.5	106	0.00
3		n-Nitrosodimethylamine	0.400	0.332	17.0	95	0.00
4	S	2-Fluorophenol	0.400	0.354	11.5	94	0.00
5	S	Phenol-d6	0.400	0.357	10.8	96	0.00
6		bis(2-Chloroethyl)ether	0.400	0.388	3.0	95	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	103	0.00
8	S	Nitrobenzene-d5	0.400	0.350	12.5	96	0.01
9		Naphthalene	0.400	0.408	-2.0	103	0.00
10		Hexachlorobutadiene	0.400	0.422	-5.5	105	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.405	-1.3	101	0.00
12		2-Methylnaphthalene	0.400	0.408	-2.0	102	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	104	0.00
14	S	2,4,6-Tribromophenol	0.400	0.335	16.3	96	0.00
15	S	2-Fluorobiphenyl	0.400	0.398	0.5	103	0.00
16		Acenaphthylene	0.400	0.380	5.0	100	0.00
17		Acenaphthene	0.400	0.407	-1.7	103	0.00
18		Fluorene	0.400	0.405	-1.3	103	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	103	0.00
20		4-Bromophenyl-phenylether	0.400	0.392	2.0	101	0.00
21		Hexachlorobenzene	0.400	0.418	-4.5	104	0.00
22		Pentachlorophenol	0.400	0.383	4.3	88	0.00
23		Phenanthrene	0.400	0.403	-0.8	102	0.01
24		Anthracene	0.400	0.374	6.5	99	0.01
25	SURR	Fluoranthene-d10	0.400	0.394	1.5	100	0.00
26		Fluoranthene	0.400	0.402	-0.5	100	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	98	-0.02
28		Pyrene	0.400	0.415	-3.7	99	0.00
29	S	Terphenyl-d14	0.400	0.410	-2.5	100	0.00
30		Benzo(a)anthracene	0.400	0.394	1.5	98	-0.02
31		Chrysene	0.400	0.397	0.8	92	-0.02
32		Bis(2-ethylhexyl)phthalate	0.400	0.322	19.5	88	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.387	3.3	81	-0.04
34	I	Perylene-d12	0.400	0.400	0.0	86	-0.04
35		Benzo(b)fluoranthene	0.400	0.427	-6.7	92	-0.03
36		Benzo(k)fluoranthene	0.400	0.427	-6.7	90	-0.04
37	C	Benzo(a)pyrene	0.400	0.419	-4.7	87	-0.04
38		Dibenzo(a,h)anthracene	0.400	0.400	0.0	81	-0.04
39		Benzo(g,h,i)perylene	0.400	0.398	0.5	81	-0.05

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

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