Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE061119\
Data File : BE099925.D
Acq On : 11 Jun 2019 15:43
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 12 04:17:38 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE061019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 10 14:57:20 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	99	-0.01
2		1,4-Dioxane	0.400	0.418	-4.5	111	0.00
3		n-Nitrosodimethylamine	0.400	0.281	29.8#	69	0.00
4	S	2-Fluorophenol	0.400	0.374	6.5	93	0.00
5	S	Phenol-d6	0.400	0.409	-2.2	109	0.00
6		bis(2-Chloroethyl)ether	0.400	0.414	-3.5	108	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	104	-0.01
8	S	Nitrobenzene-d5	0.400	0.351	12.3	97	-0.01
9		Naphthalene	0.400	0.418	-4.5	110	-0.03
10		Hexachlorobutadiene	0.400	0.420	-5.0	107	-0.03
11	SURR	2-Methylnaphthalene-d10	0.400	0.413	-3.2	110	-0.03
12		2-Methylnaphthalene	0.400	0.407	-1.7	107	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	116	-0.03
14	S	2,4,6-Tribromophenol	0.400	0.294	26.5#	94	-0.03
15	S	2-Fluorobiphenyl	0.400	0.376	6.0	115	-0.03
16		Acenaphthylene	0.400	0.387	3.3	116	-0.03
17		Acenaphthene	0.400	0.383	4.3	115	-0.03
18		Fluorene	0.400	0.409	-2.2	125	-0.03
19	I	Phenanthrene-d10	0.400	0.400	0.0	121	-0.03
20		4-Bromophenyl-phenylether	0.400	0.384	4.0	126	-0.03
21		Hexachlorobenzene	0.400	0.395	1.3	127	-0.04
22		Pentachlorophenol	0.400	0.421	-5.2	64	-0.04
23		Phenanthrene	0.400	0.384	4.0	125	-0.03
24		Anthracene	0.400	0.363	9.3	116	-0.03
25	SURR	Fluoranthene-d10	0.400	0.349	12.8	108	-0.04
26		Fluoranthene	0.400	0.356	11.0	110	-0.04
27	I	Chrysene-d12	0.400	0.400	0.0	101	-0.04
28		Pyrene	0.400	0.461	-15.3	109	-0.03
29	S	Terphenyl-d14	0.400	0.435	-8.7	105	-0.03
30		Benzo(a)anthracene	0.400	0.406	-1.5	99	-0.04
31		Chrysene	0.400	0.427	-6.7	99	-0.04
32		Bis(2-ethylhexyl)phthalate	0.400	0.376	6.0	93	-0.04
33		Indeno(1,2,3-cd)pyrene	0.400	0.411	-2.7	98	-0.10
34	I	Perylene-d12	0.400	0.400	0.0	100	-0.07
35		Benzo(b)fluoranthene	0.400	0.381	4.8	100	-0.06
36		Benzo(k)fluoranthene	0.400	0.413	-3.2	103	-0.06
37	C	Benzo(a)pyrene	0.400	0.391	2.3	101	-0.06
38		Dibenzo(a,h)anthracene	0.400	0.357	10.8	94	-0.09
39		Benzo(g,h,i)perylene	0.400	0.395	1.3	101	-0.11

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

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