Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE061119\
Data File : BE099927.D
Acq On : 11 Jun 2019 16:57
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 12 16:38:20 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	111	-0.01
2		1,4-Dioxane	0.400	0.350	12.5	103	0.00
3		n-Nitrosodimethylamine	0.400	0.358	10.5	99	0.03
4	S	2-Fluorophenol	0.400	0.347	13.3	96	0.00
5	S	Phenol-d6	0.400	0.379	5.3	113	0.00
6		bis(2-Chloroethyl)ether	0.400	0.389	2.8	114	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	110	-0.01
8	S	Nitrobenzene-d5	0.400	0.356	11.0	104	0.00
9		Naphthalene	0.400	0.411	-2.7	115	-0.03
10		Hexachlorobutadiene	0.400	0.417	-4.2	113	-0.03
11	SURR	2-Methylnaphthalene-d10	0.400	0.407	-1.7	114	-0.01
12		2-Methylnaphthalene	0.400	0.406	-1.5	113	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	120	-0.03
14	S	2,4,6-Tribromophenol	0.400	0.348	13.0	116	-0.03
15	S	2-Fluorobiphenyl	0.400	0.378	5.5	120	-0.03
16		Acenaphthylene	0.400	0.389	2.8	121	-0.03
17		Acenaphthene	0.400	0.374	6.5	117	-0.03
18		Fluorene	0.400	0.392	2.0	124	-0.04
19	I	Phenanthrene-d10	0.400	0.400	0.0	123	-0.03
20		4-Bromophenyl-phenylether	0.400	0.386	3.5	128	-0.03
21		Hexachlorobenzene	0.400	0.402	-0.5	130	-0.04
22		Pentachlorophenol	0.400	0.415	-3.7	62	-0.03
23		Phenanthrene	0.400	0.380	5.0	125	-0.03
24		Anthracene	0.400	0.350	12.5	114	-0.03
25	SURR	Fluoranthene-d10	0.400	0.348	13.0	109	-0.04
26		Fluoranthene	0.400	0.354	11.5	110	-0.04
27	I	Chrysene-d12	0.400	0.400	0.0	100	-0.04
28		Pyrene	0.400	0.467	-16.8	109	-0.04
29	S	Terphenyl-d14	0.400	0.443	-10.7	105	-0.03
30		Benzo(a)anthracene	0.400	0.412	-3.0	99	-0.04
31		Chrysene	0.400	0.423	-5.7	97	-0.04
32		Bis(2-ethylhexyl)phthalate	0.400	0.350	12.5	86	-0.04
33		Indeno(1,2,3-cd)pyrene	0.400	0.412	-3.0	98	-0.11
34	I	Perylene-d12	0.400	0.400	0.0	98	-0.07
35		Benzo(b)fluoranthene	0.400	0.380	5.0	98	-0.06
36		Benzo(k)fluoranthene	0.400	0.390	2.5	96	-0.06
37	C	Benzo(a)pyrene	0.400	0.397	0.8	101	-0.07
38		Dibenzo(a,h)anthracene	0.400	0.360	10.0	93	-0.10
39		Benzo(g,h,i)perylene	0.400	0.396	1.0	100	-0.11

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

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