Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE040219\
Data File : BE099284.D
Acq On : 2 Apr 2019 21:21
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc : LOQ 0.2 PPM
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Apr 03 05:18:31 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE040219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 02 17:16:42 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	98	-0.02
2		1,4-Dioxane	0.400	0.415	-3.7	102	-0.02
3		n-Nitrosodimethylamine	0.400	0.340	15.0	86	-0.02
4	S	2-Fluorophenol	0.400	0.380	5.0	89	-0.02
5	S	Phenol-d6	0.400	0.383	4.3	93	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.392	2.0	96	-0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	91	-0.01
8	S	Nitrobenzene-d5	0.400	0.417	-4.2	99	-0.03
9		Naphthalene	0.400	0.412	-3.0	93	-0.03
10		Hexachlorobutadiene	0.400	0.404	-1.0	91	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.397	0.8	92	-0.01
12		2-Methylnaphthalene	0.400	0.399	0.3	92	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	90	-0.03
14	S	2,4,6-Tribromophenol	0.400	0.350	12.5	88	-0.01
15	S	2-Fluorobiphenyl	0.400	0.425	-6.2	95	-0.01
16		Acenaphthylene	0.400	0.368	8.0	85	-0.01
17		Acenaphthene	0.400	0.401	-0.3	92	-0.01
18		Fluorene	0.400	0.396	1.0	88	-0.03
19	I	Phenanthrene-d10	0.400	0.400	0.0	89	-0.03
20		4-Bromophenyl-phenylether	0.400	0.378	5.5	86	-0.03
21		Hexachlorobenzene	0.400	0.411	-2.7	93	-0.03
22		Pentachlorophenol	0.400	0.288	28.0#	73	-0.03
23		Phenanthrene	0.400	0.402	-0.5	88	-0.03
24		Anthracene	0.400	0.371	7.3	83	-0.01
25	SURR	Fluoranthene-d10	0.400	0.378	5.5	83	-0.02
26		Fluoranthene	0.400	0.380	5.0	84	-0.02
27	I	Chrysene-d12	0.400	0.400	0.0	81	-0.02
28		Pyrene	0.400	0.399	0.3	82	-0.02
29	S	Terphenyl-d14	0.400	0.406	-1.5	83	-0.02
30		Benzo(a)anthracene	0.400	0.384	4.0	78	-0.02
31		Chrysene	0.400	0.405	-1.3	82	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.280	30.0#	59	-0.03
33		Indeno(1,2,3-cd)pyrene	0.400	0.403	-0.8	80	-0.06
34	I	Perylene-d12	0.400	0.400	0.0	79	-0.03
35		Benzo(b)fluoranthene	0.400	0.403	-0.8	83	-0.03
36		Benzo(k)fluoranthene	0.400	0.407	-1.7	83	-0.03
37	C	Benzo(a)pyrene	0.400	0.386	3.5	79	-0.03
38		Dibenzo(a,h)anthracene	0.400	0.401	-0.3	81	-0.06
39		Benzo(g,h,i)perylene	0.400	0.400	0.0	81	-0.06

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

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