Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE040319\
Data File : BE099286.D
Acq On : 3 Apr 2019 10:46
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Apr 04 09:16:43 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	105	-0.02
2		1,4-Dioxane	0.400	0.470	-17.5	124	-0.02
3		n-Nitrosodimethylamine	0.400	0.386	3.5	105	-0.02
4	S	2-Fluorophenol	0.400	0.422	-5.5	106	-0.02
5	S	Phenol-d6	0.400	0.408	-2.0	106	-0.01
6		bis(2-Chloroethyl)ether	0.400	0.427	-6.7	112	-0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	103	-0.01
8	S	Nitrobenzene-d5	0.400	0.436	-9.0	117	-0.01
9		Naphthalene	0.400	0.430	-7.5	110	-0.03
10		Hexachlorobutadiene	0.400	0.433	-8.2	111	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.414	-3.5	109	-0.01
12		2-Methylnaphthalene	0.400	0.426	-6.5	111	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	102	-0.03
14	S	2,4,6-Tribromophenol	0.400	0.369	7.8	105	-0.01
15	S	2-Fluorobiphenyl	0.400	0.434	-8.5	111	-0.01
16		Acenaphthylene	0.400	0.394	1.5	104	-0.01
17		Acenaphthene	0.400	0.419	-4.7	109	-0.01
18		Fluorene	0.400	0.415	-3.7	106	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	103	-0.03
20		4-Bromophenyl-phenylether	0.400	0.401	-0.3	105	-0.03
21		Hexachlorobenzene	0.400	0.426	-6.5	110	-0.03
22		Pentachlorophenol	0.400	0.473	-18.2	139	-0.03
23		Phenanthrene	0.400	0.427	-6.7	108	-0.03
24		Anthracene	0.400	0.412	-3.0	106	-0.01
25	SURR	Fluoranthene-d10	0.400	0.410	-2.5	104	-0.02
26		Fluoranthene	0.400	0.416	-4.0	106	-0.02
27	I	Chrysene-d12	0.400	0.400	0.0	98	-0.02
28		Pyrene	0.400	0.422	-5.5	105	-0.02
29	S	Terphenyl-d14	0.400	0.432	-8.0	107	-0.02
30		Benzo(a)anthracene	0.400	0.414	-3.5	101	-0.02
31		Chrysene	0.400	0.430	-7.5	104	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.321	19.8	81	-0.02
33		Indeno(1,2,3-cd)pyrene	0.400	0.405	-1.3	96	-0.06
34	I	Perylene-d12	0.400	0.400	0.0	92	-0.03
35		Benzo(b)fluoranthene	0.400	0.420	-5.0	100	-0.03
36		Benzo(k)fluoranthene	0.400	0.439	-9.7	104	-0.03
37	C	Benzo(a)pyrene	0.400	0.412	-3.0	98	-0.03
38		Dibenzo(a,h)anthracene	0.400	0.419	-4.7	99	-0.06
39		Benzo(g,h,i)perylene	0.400	0.414	-3.5	98	-0.05

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

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