Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE040519\
Data File : BE099302.D
Acq On : 5 Apr 2019 10:18
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Apr 05 14:59:55 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\8270-SIM-BE040219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 05 14:55:44 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	-0.02
2		1,4-Dioxane	0.400	0.441	-10.2	123	-0.01
3		n-Nitrosodimethylamine	0.400	0.340	15.0	98	-0.02
4	S	2-Fluorophenol	0.400	0.412	-3.0	110	-0.02
5	S	Phenol-d6	0.400	0.406	-1.5	112	-0.01
6		bis(2-Chloroethyl)ether	0.400	0.412	-3.0	114	-0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	109	-0.01
8	S	Nitrobenzene-d5	0.400	0.460	-15.0	131	-0.03
9		Naphthalene	0.400	0.426	-6.5	115	-0.03
10		Hexachlorobutadiene	0.400	0.438	-9.5	119	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.424	-6.0	118	-0.03
12		2-Methylnaphthalene	0.400	0.419	-4.7	116	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	108	-0.03
14	S	2,4,6-Tribromophenol	0.400	0.392	2.0	118	-0.01
15	S	2-Fluorobiphenyl	0.400	0.436	-9.0	118	-0.01
16		Acenaphthylene	0.400	0.396	1.0	111	-0.01
17		Acenaphthene	0.400	0.426	-6.5	118	-0.01
18		Fluorene	0.400	0.418	-4.5	112	-0.03
19	I	Phenanthrene-d10	0.400	0.400	0.0	107	-0.03
20		4-Bromophenyl-phenylether	0.400	0.421	-5.2	115	-0.03
21		Hexachlorobenzene	0.400	0.440	-10.0	119	-0.03
22		Pentachlorophenol	0.400	0.506	-26.5#	155	-0.03
23		Phenanthrene	0.400	0.445	-11.2	117	-0.03
24		Anthracene	0.400	0.407	-1.7	109	-0.01
25	SURR	Fluoranthene-d10	0.400	0.418	-4.5	110	-0.02
26		Fluoranthene	0.400	0.427	-6.7	113	-0.02
27	I	Chrysene-d12	0.400	0.400	0.0	101	-0.02
28		Pyrene	0.400	0.434	-8.5	112	-0.02
29	S	Terphenyl-d14	0.400	0.436	-9.0	111	-0.02
30		Benzo(a)anthracene	0.400	0.409	-2.2	103	-0.02
31		Chrysene	0.400	0.427	-6.7	107	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.308	23.0	81	-0.02
33		Indeno(1,2,3-cd)pyrene	0.400	0.412	-3.0	101	-0.06
34	I	Perylene-d12	0.400	0.400	0.0	97	-0.03
35		Benzo(b)fluoranthene	0.400	0.420	-5.0	105	-0.03
36		Benzo(k)fluoranthene	0.400	0.429	-7.2	106	-0.03
37	C	Benzo(a)pyrene	0.400	0.410	-2.5	102	-0.03
38		Dibenzo(a,h)anthracene	0.400	0.420	-5.0	104	-0.07
39		Benzo(g,h,i)perylene	0.400	0.414	-3.5	103	-0.07

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

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