Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE050619\
Data File : BE099555.D
Acq On : 6 May 2019 17:50
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 07 08:33:22 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE042919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue May 07 08:07:29 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	165	0.01
2		1,4-Dioxane	0.400	0.350	12.5	168	0.00
3		n-Nitrosodimethylamine	0.400	0.520	-30.0#	242	-0.01
4	S	2-Fluorophenol	0.400	0.401	-0.3	173	0.00
5	S	Phenol-d6	0.400	0.437	-9.2	184	0.01
6		bis(2-Chloroethyl)ether	0.400	0.398	0.5	168	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	172	0.03
8	S	Nitrobenzene-d5	0.400	0.446	-11.5	191	0.01
9		Naphthalene	0.400	0.404	-1.0	170	0.03
10		Hexachlorobutadiene	0.400	0.415	-3.7	174	0.03
11	SURR	2-Methylnaphthalene-d10	0.400	0.397	0.8	172	0.03
12		2-Methylnaphthalene	0.400	0.398	0.5	171	0.04
13	I	Acenaphthene-d10	0.400	0.400	0.0	163	0.04
14	S	2,4,6-Tribromophenol	0.400	0.487	-21.7	205	0.04
15	S	2-Fluorobiphenyl	0.400	0.416	-4.0	166	0.03
16		Acenaphthylene	0.400	0.418	-4.5	170	0.03
17		Acenaphthene	0.400	0.412	-3.0	168	0.04
18		Fluorene	0.400	0.407	-1.7	164	0.05
19	I	Phenanthrene-d10	0.400	0.400	0.0	162	0.05
20		4-Bromophenyl-phenylether	0.400	0.406	-1.5	162	0.05
21		Hexachlorobenzene	0.400	0.416	-4.0	167	0.05
22		Pentachlorophenol	0.400	0.564	-41.0#	223	0.06
23		Phenanthrene	0.400	0.397	0.8	158	0.05
24		Anthracene	0.400	0.401	-0.3	163	0.05
25	SURR	Fluoranthene-d10	0.400	0.404	-1.0	163	0.06
26		Fluoranthene	0.400	0.410	-2.5	164	0.05
27	I	Chrysene-d12	0.400	0.400	0.0	155	0.06
28		Pyrene	0.400	0.474	-18.5	162	0.06
29	S	Terphenyl-d14	0.400	0.470	-17.5	162	0.06
30		Benzo(a)anthracene	0.400	0.449	-12.2	154	0.06
31		Chrysene	0.400	0.451	-12.7	154	0.06
32		Bis(2-ethylhexyl)phthalate	0.400	0.451	-12.7	151	0.06
33		Indeno(1,2,3-cd)pyrene	0.400	0.459	-14.8	156	0.17
34	I	Perylene-d12	0.400	0.400	0.0	152	0.12
35		Benzo(b)fluoranthene	0.400	0.418	-4.5	156	0.10
36		Benzo(k)fluoranthene	0.400	0.412	-3.0	149	0.10
37	C	Benzo(a)pyrene	0.400	0.421	-5.2	155	0.11
38		Dibenzo(a,h)anthracene	0.400	0.430	-7.5	158	0.18
39		Benzo(g,h,i)perylene	0.400	0.427	-6.7	156	0.20

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

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