Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE050719\
Data File : BE099571.D
Acq On : 7 May 2019 19:14
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: May 08 02:00:01 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE050719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue May 07 14:10:11 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	121	0.00
2		1,4-Dioxane	0.400	0.390	2.5	143	0.00
3		n-Nitrosodimethylamine	0.400	0.365	8.8	122	0.00
4	S	2-Fluorophenol	0.400	0.424	-6.0	143	0.00
5	S	Phenol-d6	0.400	0.426	-6.5	141	0.00
6		bis(2-Chloroethyl)ether	0.400	0.415	-3.7	141	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	121	0.00
8	S	Nitrobenzene-d5	0.400	0.415	-3.7	140	0.00
9		Naphthalene	0.400	1.969	-392.3#	658	0.00
10		Hexachlorobutadiene	0.400	0.403	-0.8	135	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.395	1.3	133	0.00
12		2-Methylnaphthalene	0.400	1.765	-341.2#	584	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	114	0.00
14	S	2,4,6-Tribromophenol	0.400	0.391	2.3	129	0.00
15	S	2-Fluorobiphenyl	0.400	0.416	-4.0	136	0.00
16		Acenaphthylene	0.400	0.672	-68.0#	216	0.00
17		Acenaphthene	0.400	0.463	-15.8	147	0.00
18		Fluorene	0.400	0.705	-76.2#	223	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	115	0.00
20		4-Bromophenyl-phenylether	0.400	0.401	-0.3	131	0.00
21		Hexachlorobenzene	0.400	0.407	-1.7	129	0.00
22		Pentachlorophenol	0.400	0.365	8.8	131	0.00
23		Phenanthrene	0.400	1.214	-203.5#	388	0.00
24		Anthracene	0.400	0.563	-40.7#	185	0.00
25	SURR	Fluoranthene-d10	0.400	0.404	-1.0	131	0.00
26		Fluoranthene	0.400	0.658	-64.5#	211	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	116	0.00
28		Pyrene	0.400	0.833	-108.2#	244	0.00
29	S	Terphenyl-d14	0.400	0.464	-16.0	136	0.00
30		Benzo(a)anthracene	0.400	0.584	-46.0#	172	-0.01
31		Chrysene	0.400	0.553	-38.3#	160	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.473	-18.2	140	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.465	-16.3	136	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	112	0.00
35		Benzo(b)fluoranthene	0.400	0.505	-26.2#	155	0.00
36		Benzo(k)fluoranthene	0.400	0.479	-19.7	148	0.00
37	C	Benzo(a)pyrene	0.400	0.512	-28.0#	157	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.431	-7.7	133	0.00
39		Benzo(g,h,i)perylene	0.400	0.452	-13.0	137	0.00

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

(#) = Out of Range SPCC's out = 0 CCC's out = 1