Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE032319\
Data File : BE099120.D
Acq On : 23 Mar 2019 20:08
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Mar 23 22:24:19 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE032219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Mar 23 04:40:11 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	99	0.00
2		1,4-Dioxane	0.400	0.423	-5.7	121	-0.01
3		n-Nitrosodimethylamine	0.400	0.446	-11.5	120	-0.01
4	S	2-Fluorophenol	0.400	0.477	-19.2	133	0.00
5	S	Phenol-d6	0.400	0.495	-23.7	141	0.00
6		bis(2-Chloroethyl)ether	0.400	0.465	-16.3	129	-0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	109	-0.01
8	S	Nitrobenzene-d5	0.400	0.837	-109.2#	286	-0.01
9		Naphthalene	0.400	0.442	-10.5	136	0.00
10		Hexachlorobutadiene	0.400	0.463	-15.8	140	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.472	-18.0	147	-0.02
12		2-Methylnaphthalene	0.400	0.472	-18.0	149	-0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	127	0.00
14	S	2,4,6-Tribromophenol	0.400	0.804	-101.0#	329	0.00
15	S	2-Fluorobiphenyl	0.400	0.455	-13.7	158	-0.02
16		Acenaphthylene	0.400	0.432	-8.0	162	-0.02
17		Acenaphthene	0.400	0.436	-9.0	162	0.00
18		Fluorene	0.400	0.440	-10.0	162	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	124	0.00
20		4-Bromophenyl-phenylether	0.400	0.485	-21.2	176	0.00
21		Hexachlorobenzene	0.400	0.489	-22.2	180	0.00
22		Pentachlorophenol	0.400	0.714	-78.5#	289	-0.01
23		Phenanthrene	0.400	0.434	-8.5	159	0.00
24		Anthracene	0.400	0.419	-4.7	156	0.00
25	SURR	Fluoranthene-d10	0.400	0.404	-1.0	144	0.00
26		Fluoranthene	0.400	0.394	1.5	146	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	116	-0.01
28		Pyrene	0.400	0.425	-6.2	146	0.00
29	S	Terphenyl-d14	0.400	0.472	-18.0	156	0.00
30		Benzo(a)anthracene	0.400	0.421	-5.2	144	0.00
31		Chrysene	0.400	0.426	-6.5	148	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.489	-22.2	201	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.438	-9.5	151	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	124	0.00
35		Benzo(b)fluoranthene	0.400	0.400	0.0	150	0.00
36		Benzo(k)fluoranthene	0.400	0.414	-3.5	154	0.00
37	C	Benzo(a)pyrene	0.400	0.404	-1.0	152	0.00
38		Dibenzo(a,h)anthracene	0.400	0.408	-2.0	154	-0.02
39		Benzo(g,h,i)perylene	0.400	0.401	-0.3	149	-0.02

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

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