Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE032319\
Data File : BE099103.D
Acq On : 23 Mar 2019 9:30
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Mar 23 22:34:47 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	91	0.00
2		1,4-Dioxane	0.400	0.494	-23.5	130	-0.01
3		n-Nitrosodimethylamine	0.400	0.369	7.8	91	-0.01
4	S	2-Fluorophenol	0.400	0.467	-16.8	120	0.00
5	S	Phenol-d6	0.400	0.487	-21.7	127	0.00
6		bis(2-Chloroethyl)ether	0.400	0.472	-18.0	121	-0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	101	0.00
8	S	Nitrobenzene-d5	0.400	0.722	-80.5#	230	-0.01
9		Naphthalene	0.400	0.441	-10.2	126	0.00
10		Hexachlorobutadiene	0.400	0.454	-13.5	128	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.457	-14.2	132	-0.02
12		2-Methylnaphthalene	0.400	0.463	-15.8	136	-0.02
13	I	Acenaphthene-d10	0.400	0.400	0.0	120	0.00
14	S	2,4,6-Tribromophenol	0.400	0.640	-60.0#	247	0.00
15	S	2-Fluorobiphenyl	0.400	0.435	-8.7	142	-0.02
16		Acenaphthylene	0.400	0.405	-1.3	143	-0.02
17		Acenaphthene	0.400	0.426	-6.5	149	0.00
18		Fluorene	0.400	0.424	-6.0	148	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	115	0.00
20		4-Bromophenyl-phenylether	0.400	0.457	-14.2	155	0.00
21		Hexachlorobenzene	0.400	0.471	-17.7	162	0.00
22		Pentachlorophenol	0.400	0.601	-50.2#	203	0.00
23		Phenanthrene	0.400	0.429	-7.2	147	0.00
24		Anthracene	0.400	0.405	-1.3	141	0.00
25	SURR	Fluoranthene-d10	0.400	0.373	6.8	124	0.00
26		Fluoranthene	0.400	0.371	7.3	129	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	101	-0.01
28		Pyrene	0.400	0.426	-6.5	127	0.00
29	S	Terphenyl-d14	0.400	0.471	-17.7	135	0.00
30		Benzo(a)anthracene	0.400	0.404	-1.0	120	0.00
31		Chrysene	0.400	0.428	-7.0	128	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.352	12.0	125	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.418	-4.5	125	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	100	0.00
35		Benzo(b)fluoranthene	0.400	0.416	-4.0	126	0.00
36		Benzo(k)fluoranthene	0.400	0.411	-2.7	123	0.00
37	C	Benzo(a)pyrene	0.400	0.405	-1.3	123	0.00
38		Dibenzo(a,h)anthracene	0.400	0.416	-4.0	127	-0.01
39		Benzo(g,h,i)perylene	0.400	0.418	-4.5	126	-0.01

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

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