Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA E\Data\BE041519\
Data File : BE099322.D
Acq On : 15 Apr 2019 9:33
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Apr 15 16:48:46 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE040219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 02 17:16:42 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	143	-0.02
2		1,4-Dioxane	0.400	0.415	-3.7	149	-0.01
3		n-Nitrosodimethylamine	0.400	0.419	-4.7	155	-0.02
4	S	2-Fluorophenol	0.400	0.404	-1.0	138	-0.02
5	S	Phenol-d6	0.400	0.431	-7.7	152	-0.01
6		bis(2-Chloroethyl)ether	0.400	0.430	-7.5	153	-0.02
7	I	Naphthalene-d8	0.400	0.400	0.0	145	-0.01
8	S	Nitrobenzene-d5	0.400	0.458	-14.5	173	-0.03
9		Naphthalene	0.400	0.425	-6.2	153	-0.03
10		Hexachlorobutadiene	0.400	0.412	-3.0	149	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.422	-5.5	156	-0.01
12		2-Methylnaphthalene	0.400	0.430	-7.5	157	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	152	-0.03
14	S	2,4,6-Tribromophenol	0.400	0.350	12.5	148	-0.01
15	S	2-Fluorobiphenyl	0.400	0.418	-4.5	158	-0.01
16		Acenaphthylene	0.400	0.390	2.5	153	-0.01
17		Acenaphthene	0.400	0.418	-4.5	162	-0.01
18		Fluorene	0.400	0.417	-4.2	157	-0.03
19	I	Phenanthrene-d10	0.400	0.400	0.0	152	-0.03
20		4-Bromophenyl-phenylether	0.400	0.402	-0.5	156	-0.03
21		Hexachlorobenzene	0.400	0.399	0.3	153	-0.03
22		Pentachlorophenol	0.400	0.451	-12.7	196	-0.03
23		Phenanthrene	0.400	0.429	-7.2	160	-0.03
24		Anthracene	0.400	0.412	-3.0	157	-0.01
25	SURR	Fluoranthene-d10	0.400	0.398	0.5	149	-0.02
26		Fluoranthene	0.400	0.405	-1.3	153	-0.02
27	I	Chrysene-d12	0.400	0.400	0.0	125	-0.02
28		Pyrene	0.400	0.476	-19.0	151	-0.02
29	S	Terphenyl-d14	0.400	0.458	-14.5	144	-0.02
30		Benzo(a)anthracene	0.400	0.413	-3.2	129	-0.02
31		Chrysene	0.400	0.435	-8.7	135	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.334	16.5	108	-0.02
33		Indeno(1,2,3-cd)pyrene	0.400	0.403	-0.8	122	-0.05
34	I	Perylene-d12	0.400	0.400	0.0	119	-0.03
35		Benzo(b)fluoranthene	0.400	0.416	-4.0	128	-0.03
36		Benzo(k)fluoranthene	0.400	0.437	-9.2	133	-0.03
37	C	Benzo(a)pyrene	0.400	0.414	-3.5	127	-0.03
38		Dibenzo(a,h)anthracene	0.400	0.416	-4.0	126	-0.06
39		Benzo(g,h,i)perylene	0.400	0.408	-2.0	124	-0.06

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

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