Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE121518\
Data File : BE098502.D
Acq On : 16 Dec 2018 11:08
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 34 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 17 02:05:10 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE121018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Dec 10 14:59:55 2018
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	51	-0.01
2		1,4-Dioxane	0.400	0.353	11.8	45	0.00
3		n-Nitrosodimethylamine	0.400	0.301	24.8	41	0.00
4	S	2-Fluorophenol	0.400	0.497	-24.2	65	0.01
5	S	Phenol-d6	0.400	0.493	-23.2	62	0.00
6		bis(2-Chloroethyl)ether	0.400	0.372	7.0	49	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	51	-0.01
8	S	Nitrobenzene-d5	0.400	0.394	1.5	55	0.00
9		Naphthalene	0.400	0.408	-2.0	54	0.00
10		Hexachlorobutadiene	0.400	0.470	-17.5	62	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.402	-0.5	52	-0.01
12		2-Methylnaphthalene	0.400	0.396	1.0	52	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	49	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.509	-27.2#	66	0.00
15	S	2-Fluorobiphenyl	0.400	0.410	-2.5	56	0.00
16		Acenaphthylene	0.400	0.411	-2.7	53	0.00
17		Acenaphthene	0.400	0.408	-2.0	51	0.00
18		Fluorene	0.400	0.415	-3.7	52	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	47	-0.01
20		4-Bromophenyl-phenylether	0.400	0.399	0.3	49	0.00
21		Hexachlorobenzene	0.400	0.435	-8.7	52	-0.01
22		Pentachlorophenol	0.400	0.571	-42.7#	73	0.00
23		Phenanthrene	0.400	0.415	-3.7	49	0.00
24		Anthracene	0.400	0.380	5.0	45	0.00
25	SURR	Fluoranthene-d10	0.400	0.415	-3.7	47	0.00
26		Fluoranthene	0.400	0.420	-5.0	47	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	54	-0.01
28		Pyrene	0.400	0.390	2.5	51	0.00
29	S	Terphenyl-d14	0.400	0.342	14.5	46	0.00
30		Benzo(a)anthracene	0.400	0.387	3.3	53	-0.01
31		Chrysene	0.400	0.422	-5.5	57	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.307	23.3	41	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.431	-7.7	61	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	56	-0.01
35		Benzo(b)fluoranthene	0.400	0.422	-5.5	61	0.00
36		Benzo(k)fluoranthene	0.400	0.427	-6.7	63	-0.01
37	C	Benzo(a)pyrene	0.400	0.426	-6.5	61	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.405	-1.3	59	-0.02
39		Benzo(g,h,i)perylene	0.400	0.439	-9.7	64	-0.02

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

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