Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE010419\
Data File : BE098614.D
Acq On : 4 Jan 2019 12:23
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jan 04 21:12:25 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE121018.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jan 02 10:30:27 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	64	0.00
2		1,4-Dioxane	0.400	0.404	-1.0	66	0.00
3		n-Nitrosodimethylamine	0.400	0.410	-2.5	71	0.01
4	S	2-Fluorophenol	0.400	0.531	-32.8#	88	0.01
5	S	Phenol-d6	0.400	0.559	-39.8#	89	0.01
6		bis(2-Chloroethyl)ether	0.400	0.402	-0.5	67	0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	66	0.00
8	S	Nitrobenzene-d5	0.400	0.425	-6.2	76	0.00
9		Naphthalene	0.400	0.425	-6.2	72	0.00
10		Hexachlorobutadiene	0.400	0.463	-15.8	79	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.445	-11.2	74	0.00
12		2-Methylnaphthalene	0.400	0.439	-9.7	74	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	67	0.00
14	S	2,4,6-Tribromophenol	0.400	0.507	-26.7#	90	0.01
15	S	2-Fluorobiphenyl	0.400	0.393	1.8	72	0.02
16		Acenaphthylene	0.400	0.397	0.8	69	0.02
17		Acenaphthene	0.400	0.417	-4.2	71	0.02
18		Fluorene	0.400	0.402	-0.5	68	0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	56	0.00
20		4-Bromophenyl-phenylether	0.400	0.429	-7.2	63	0.01
21		Hexachlorobenzene	0.400	0.460	-15.0	66	0.01
22		Pentachlorophenol	0.400	0.935	-133.8#	160	0.01
23		Phenanthrene	0.400	0.422	-5.5	60	0.01
24		Anthracene	0.400	0.411	-2.7	59	0.01
25	SURR	Fluoranthene-d10	0.400	0.407	-1.7	55	0.01
26		Fluoranthene	0.400	0.397	0.8	54	0.01
27	I	Chrysene-d12	0.400	0.400	0.0	61	0.01
28		Pyrene	0.400	0.381	4.8	56	0.01
29	S	Terphenyl-d14	0.400	0.386	3.5	59	0.01
30		Benzo(a)anthracene	0.400	0.425	-6.2	66	0.00
31		Chrysene	0.400	0.406	-1.5	62	0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.527	-31.8#	80	0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.433	-8.2	69	-0.06
34	I	Perylene-d12	0.400	0.400	0.0	66	-0.03
35		Benzo(b)fluoranthene	0.400	0.420	-5.0	71	-0.01
36		Benzo(k)fluoranthene	0.400	0.389	2.8	67	-0.01
37	C	Benzo(a)pyrene	0.400	0.399	0.3	67	-0.03
38		Dibenzo(a,h)anthracene	0.400	0.395	1.3	67	-0.07
39		Benzo(g,h,i)perylene	0.400	0.416	-4.0	71	-0.08

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

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