Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE121618\
Data File : BE098504.D
Acq On : 16 Dec 2018 13:14
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 17 03:16:01 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	47	0.00
2		1,4-Dioxane	0.400	0.452	-13.0	54	0.00
3		n-Nitrosodimethylamine	0.400	0.453	-13.2	58	0.00
4	S	2-Fluorophenol	0.400	0.511	-27.7#	62	0.00
5	S	Phenol-d6	0.400	0.461	-15.3	54	0.00
6		bis(2-Chloroethyl)ether	0.400	0.389	2.8	47	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	45	-0.01
8	S	Nitrobenzene-d5	0.400	0.376	6.0	46	0.00
9		Naphthalene	0.400	0.403	-0.8	46	0.00
10		Hexachlorobutadiene	0.400	0.450	-12.5	52	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.413	-3.2	47	-0.01
12		2-Methylnaphthalene	0.400	0.431	-7.7	50	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	44	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.600	-50.0#	70	-0.01
15	S	2-Fluorobiphenyl	0.400	0.413	-3.2	50	0.00
16		Acenaphthylene	0.400	0.422	-5.5	48	0.00
17		Acenaphthene	0.400	0.422	-5.5	47	0.00
18		Fluorene	0.400	0.434	-8.5	48	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	46	-0.01
20		4-Bromophenyl-phenylether	0.400	0.414	-3.5	49	-0.01
21		Hexachlorobenzene	0.400	0.429	-7.2	50	-0.01
22		Pentachlorophenol	0.400	1.015	-153.7#	144	-0.01
23		Phenanthrene	0.400	0.419	-4.7	49	0.00
24		Anthracene	0.400	0.409	-2.2	47	0.00
25	SURR	Fluoranthene-d10	0.400	0.481	-20.2	53	0.00
26		Fluoranthene	0.400	0.477	-19.2	52	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	60	0.00
28		Pyrene	0.400	0.387	3.3	57	0.00
29	S	Terphenyl-d14	0.400	0.363	9.3	55	0.00
30		Benzo(a)anthracene	0.400	0.436	-9.0	67	-0.01
31		Chrysene	0.400	0.408	-2.0	62	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.556	-39.0#	84	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.412	-3.0	65	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	63	-0.01
35		Benzo(b)fluoranthene	0.400	0.436	-9.0	70	-0.01
36		Benzo(k)fluoranthene	0.400	0.406	-1.5	66	-0.01
37	C	Benzo(a)pyrene	0.400	0.424	-6.0	68	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.385	3.8	62	-0.01
39		Benzo(g,h,i)perylene	0.400	0.417	-4.2	67	-0.02

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

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