Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE121518\
Data File : BE098485.D
Acq On : 15 Dec 2018 23:09
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 16 02:58:19 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	46	0.00
2		1,4-Dioxane	0.400	0.348	13.0	41	0.00
3		n-Nitrosodimethylamine	0.400	0.221	44.8#	28	0.00
4	S	2-Fluorophenol	0.400	0.453	-13.2	54	0.00
5	S	Phenol-d6	0.400	0.416	-4.0	48	0.00
6		bis(2-Chloroethyl)ether	0.400	0.351	12.3	42	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	45	-0.01
8	S	Nitrobenzene-d5	0.400	0.363	9.3	44	0.01
9		Naphthalene	0.400	0.415	-3.7	48	0.00
10		Hexachlorobutadiene	0.400	0.467	-16.8	54	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.406	-1.5	46	0.00
12		2-Methylnaphthalene	0.400	0.389	2.8	45	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	44	0.00
14	S	2,4,6-Tribromophenol	0.400	0.396	1.0	46	0.00
15	S	2-Fluorobiphenyl	0.400	0.403	-0.8	49	0.00
16		Acenaphthylene	0.400	0.419	-4.7	48	0.00
17		Acenaphthene	0.400	0.404	-1.0	45	0.00
18		Fluorene	0.400	0.416	-4.0	47	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	42	-0.01
20		4-Bromophenyl-phenylether	0.400	0.373	6.8	41	0.00
21		Hexachlorobenzene	0.400	0.437	-9.2	47	0.00
22		Pentachlorophenol	0.400	0.368	8.0	39	0.01
23		Phenanthrene	0.400	0.403	-0.8	43	0.00
24		Anthracene	0.400	0.343	14.2	37	0.00
25	SURR	Fluoranthene-d10	0.400	0.410	-2.5	42	0.00
26		Fluoranthene	0.400	0.415	-3.7	42	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	46	0.00
28		Pyrene	0.400	0.375	6.3	42	0.00
29	S	Terphenyl-d14	0.400	0.332	17.0	38	0.00
30		Benzo(a)anthracene	0.400	0.376	6.0	44	-0.01
31		Chrysene	0.400	0.400	0.0	46	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.262	34.5#	30	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.372	7.0	45	0.00
34	I	Perylene-d12	0.400	0.400	0.0	48	-0.01
35		Benzo(b)fluoranthene	0.400	0.394	1.5	49	0.00
36		Benzo(k)fluoranthene	0.400	0.443	-10.7	55	0.00
37	C	Benzo(a)pyrene	0.400	0.406	-1.5	50	0.00
38		Dibenzo(a,h)anthracene	0.400	0.345	13.8	43	0.00
39		Benzo(g,h,i)perylene	0.400	0.409	-2.2	51	0.00

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

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