Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE121518\
Data File : BE098468.D
Acq On : 15 Dec 2018 12:29
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Dec 16 02:33:09 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	53	0.00
2		1,4-Dioxane	0.400	0.343	14.2	46	0.00
3		n-Nitrosodimethylamine	0.400	0.277	30.8#	39	0.01
4	S	2-Fluorophenol	0.400	0.478	-19.5	65	0.01
5	S	Phenol-d6	0.400	0.495	-23.7	65	0.00
6		bis(2-Chloroethyl)ether	0.400	0.375	6.3	51	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	53	-0.01
8	S	Nitrobenzene-d5	0.400	0.382	4.5	56	0.01
9		Naphthalene	0.400	0.411	-2.7	56	0.00
10		Hexachlorobutadiene	0.400	0.450	-12.5	62	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.407	-1.7	55	0.00
12		2-Methylnaphthalene	0.400	0.382	4.5	52	0.00
13	I	Acenaphthene-d10	0.400	0.400	0.0	52	0.00
14	S	2,4,6-Tribromophenol	0.400	0.471	-17.7	65	0.00
15	S	2-Fluorobiphenyl	0.400	0.385	3.8	55	0.00
16		Acenaphthylene	0.400	0.417	-4.2	56	0.00
17		Acenaphthene	0.400	0.408	-2.0	54	0.00
18		Fluorene	0.400	0.408	-2.0	54	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	49	-0.01
20		4-Bromophenyl-phenylether	0.400	0.380	5.0	49	0.00
21		Hexachlorobenzene	0.400	0.422	-5.5	53	0.00
22		Pentachlorophenol	0.400	0.543	-35.8#	72	0.00
23		Phenanthrene	0.400	0.408	-2.0	51	0.00
24		Anthracene	0.400	0.372	7.0	46	0.00
25	SURR	Fluoranthene-d10	0.400	0.423	-5.7	50	0.00
26		Fluoranthene	0.400	0.425	-6.2	50	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	57	0.00
28		Pyrene	0.400	0.381	4.8	53	0.00
29	S	Terphenyl-d14	0.400	0.335	16.3	48	0.00
30		Benzo(a)anthracene	0.400	0.403	-0.8	59	-0.01
31		Chrysene	0.400	0.407	-1.7	59	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.388	3.0	56	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.408	-2.0	61	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	61	-0.01
35		Benzo(b)fluoranthene	0.400	0.404	-1.0	63	0.00
36		Benzo(k)fluoranthene	0.400	0.429	-7.2	68	-0.01
37	C	Benzo(a)pyrene	0.400	0.396	1.0	61	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.384	4.0	60	0.00
39		Benzo(g,h,i)perylene	0.400	0.401	-0.3	63	-0.01

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

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