Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE101618\
Data File : BE097930.D
Acq On : 17 Oct 2018 3:28
Operator : SJ/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Oct 17 03:59:01 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE101618.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 16 15:36:44 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	111	0.00
2		1,4-Dioxane	0.400	0.436	-9.0	131	0.00
3		n-Nitrosodimethylamine	0.400	0.371	7.3	117	0.00
4	S	2-Fluorophenol	0.400	0.413	-3.2	122	-0.01
5	S	Phenol-d6	0.400	0.422	-5.5	121	-0.01
6		bis(2-Chloroethyl)ether	0.400	0.426	-6.5	127	0.00
7	I	Naphthalene-d8	0.400	0.400	0.0	117	-0.01
8	S	Nitrobenzene-d5	0.400	0.415	-3.7	124	-0.01
9		Naphthalene	0.400	0.437	-9.2	131	-0.01
10		Hexachlorobutadiene	0.400	0.427	-6.7	129	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.440	-10.0	132	-0.01
12		2-Methylnaphthalene	0.400	0.438	-9.5	133	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	124	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.408	-2.0	143	0.00
15	S	2-Fluorobiphenyl	0.400	0.406	-1.5	133	0.00
16		Acenaphthylene	0.400	0.430	-7.5	141	0.00
17		Acenaphthene	0.400	0.429	-7.2	140	0.00
18		Fluorene	0.400	0.437	-9.2	142	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	138	-0.01
20		4-Bromophenyl-phenylether	0.400	0.420	-5.0	148	0.00
21		Hexachlorobenzene	0.400	0.428	-7.0	151	0.00
22		Pentachlorophenol	0.400	0.452	-13.0	176	-0.01
23		Phenanthrene	0.400	0.427	-6.7	152	0.00
24		Anthracene	0.400	0.421	-5.2	151	0.00
25	SURR	Fluoranthene-d10	0.400	0.426	-6.5	158	0.00
26		Fluoranthene	0.400	0.430	-7.5	159	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	120	-0.01
28		Pyrene	0.400	0.494	-23.5	157	0.00
29	S	Terphenyl-d14	0.400	0.475	-18.7	157	0.00
30		Benzo(a)anthracene	0.400	0.422	-5.5	134	0.00
31		Chrysene	0.400	0.437	-9.2	143	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.293	26.8#	99	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.335	16.3	110	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	99	-0.01
35		Benzo(b)fluoranthene	0.400	0.420	-5.0	111	-0.01
36		Benzo(k)fluoranthene	0.400	0.448	-12.0	122	0.00
37	C	Benzo(a)pyrene	0.400	0.425	-6.2	111	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.427	-6.7	112	-0.01
39		Benzo(g,h,i)perylene	0.400	0.414	-3.5	112	-0.01

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

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