Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE032319\
Data File : BE099139.D
Acq On : 24 Mar 2019 8:40
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Mar 24 23:03:18 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE032219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Mar 23 04:40:11 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	148	0.00
2		1,4-Dioxane	0.400	0.388	3.0	165	-0.01
3		n-Nitrosodimethylamine	0.400	0.460	-15.0	184	-0.02
4	S	2-Fluorophenol	0.400	0.468	-17.0	194	0.00
5	S	Phenol-d6	0.400	0.502	-25.5#	212	0.00
6		bis(2-Chloroethyl)ether	0.400	0.468	-17.0	193	-0.01
7	I	Naphthalene-d8	0.400	0.400	0.0	170	-0.01
8	S	Nitrobenzene-d5	0.400	0.906	-126.5#	483	0.00
9		Naphthalene	0.400	0.439	-9.7	211	0.00
10		Hexachlorobutadiene	0.400	0.479	-19.7	226	0.00
11	SURR	2-Methylnaphthalene-d10	0.400	0.464	-16.0	225	-0.01
12		2-Methylnaphthalene	0.400	0.464	-16.0	228	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	187	0.00
14	S	2,4,6-Tribromophenol	0.400	0.851	-112.7#	512	0.00
15	S	2-Fluorobiphenyl	0.400	0.468	-17.0	238	-0.01
16		Acenaphthylene	0.400	0.436	-9.0	241	-0.01
17		Acenaphthene	0.400	0.428	-7.0	234	0.00
18		Fluorene	0.400	0.410	-2.5	222	0.00
19	I	Phenanthrene-d10	0.400	0.400	0.0	170	0.00
20		4-Bromophenyl-phenylether	0.400	0.480	-20.0	240	0.00
21		Hexachlorobenzene	0.400	0.487	-21.7	246	0.00
22		Pentachlorophenol	0.400	0.749	-87.3#	426	-0.01
23		Phenanthrene	0.400	0.416	-4.0	209	0.00
24		Anthracene	0.400	0.414	-3.5	212	-0.01
25	SURR	Fluoranthene-d10	0.400	0.391	2.3	192	0.00
26		Fluoranthene	0.400	0.382	4.5	195	0.00
27	I	Chrysene-d12	0.400	0.400	0.0	156	0.00
28		Pyrene	0.400	0.425	-6.2	196	0.00
29	S	Terphenyl-d14	0.400	0.479	-19.7	212	0.00
30		Benzo(a)anthracene	0.400	0.422	-5.5	194	0.00
31		Chrysene	0.400	0.428	-7.0	198	0.00
32		Bis(2-ethylhexyl)phthalate	0.400	0.571	-42.7#	314	-0.01
33		Indeno(1,2,3-cd)pyrene	0.400	0.447	-11.7	207	-0.01
34	I	Perylene-d12	0.400	0.400	0.0	166	0.00
35		Benzo(b)fluoranthene	0.400	0.403	-0.8	203	0.00
36		Benzo(k)fluoranthene	0.400	0.405	-1.3	201	-0.01
37	C	Benzo(a)pyrene	0.400	0.409	-2.2	207	0.00
38		Dibenzo(a,h)anthracene	0.400	0.419	-4.7	212	-0.02
39		Benzo(g,h,i)perylene	0.400	0.408	-2.0	203	-0.01

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

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