Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE062119\
Data File : BE100154.D
Acq On : 22 Jun 2019 4:00
Operator : JU/SJ
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jun 22 04:40:55 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE061919.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 19 15:07:24 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	169	-0.01
2		1,4-Dioxane	0.400	0.255	36.3#	122	-0.01
3		n-Nitrosodimethylamine	0.400	0.593	-48.2#	638	-0.06
4	S	2-Fluorophenol	0.400	0.535	-33.8#	239	-0.01
5	S	Phenol-d6	0.400	0.587	-46.7#	245	-0.02
6		bis(2-Chloroethyl)ether	0.400	0.447	-11.7	205	-0.03
7	I	Naphthalene-d8	0.400	0.400	0.0	190	-0.01
8	S	Nitrobenzene-d5	0.400	0.435	-8.7	205	-0.02
9		Naphthalene	0.400	0.406	-1.5	194	-0.01
10		Hexachlorobutadiene	0.400	0.367	8.3	179	-0.01
11	SURR	2-Methylnaphthalene-d10	0.400	0.407	-1.7	202	-0.01
12		2-Methylnaphthalene	0.400	0.431	-7.7	214	-0.01
13	I	Acenaphthene-d10	0.400	0.400	0.0	202	-0.01
14	S	2,4,6-Tribromophenol	0.400	0.525	-31.3#	297	-0.01
15	S	2-Fluorobiphenyl	0.400	0.417	-4.2	194	-0.03
16		Acenaphthylene	0.400	0.434	-8.5	217	-0.01
17		Acenaphthene	0.400	0.419	-4.7	207	0.00
18		Fluorene	0.400	0.401	-0.3	199	-0.01
19	I	Phenanthrene-d10	0.400	0.400	0.0	175	-0.01
20		4-Bromophenyl-phenylether	0.400	0.441	-10.2	200	-0.01
21		Hexachlorobenzene	0.400	0.415	-3.7	182	-0.01
22		Pentachlorophenol	0.400	0.724	-81.0#	466	-0.05
23		Phenanthrene	0.400	0.411	-2.7	182	-0.01
24		Anthracene	0.400	0.447	-11.7	210	-0.01
25	SURR	Fluoranthene-d10	0.400	0.368	8.0	163	0.00
26		Fluoranthene	0.400	0.366	8.5	162	-0.01
27	I	Chrysene-d12	0.400	0.400	0.0	136	0.00
28		Pyrene	0.400	0.492	-23.0	169	-0.01
29	S	Terphenyl-d14	0.400	0.495	-23.7	168	-0.02
30		Benzo(a)anthracene	0.400	0.455	-13.7	169	0.00
31		Chrysene	0.400	0.393	1.8	130	-0.01
32		Bis(2-ethylhexyl)phthalate	0.400	0.813	-103.2#	320	0.00
33		Indeno(1,2,3-cd)pyrene	0.400	0.419	-4.7	137	0.00
34	I	Perylene-d12	0.400	0.400	0.0	147	-0.01
35		Benzo(b)fluoranthene	0.400	0.373	6.8	138	0.00
36		Benzo(k)fluoranthene	0.400	0.345	13.8	131	0.00
37	C	Benzo(a)pyrene	0.400	0.373	6.8	137	-0.01
38		Dibenzo(a,h)anthracene	0.400	0.358	10.5	136	0.00
39		Benzo(g,h,i)perylene	0.400	0.354	11.5	131	0.00

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

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