Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE072319\
Data File : BE100491.D
Acq On : 23 Jul 2019 8:28
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC0.4

Quant Time: Jul 24 02:20:17 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\8270-SIM-BE072219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 23 08:06:59 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	99	0.00
2		1,4-Dioxane	0.601	0.550	8.5	91	0.00
3		n-Nitrosodimethylamine	0.366	0.312	14.8	84	0.00
4	S	2-Fluorophenol	0.976	0.948	2.9	97	0.00
5	S	Phenol-d6	1.394	1.369	1.8	98	0.00
6		bis(2-Chloroethyl)ether	1.196	1.218	-1.8	102	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	101	0.00
8	S	Nitrobenzene-d5	0.332	0.308	7.2	99	0.00
9		Naphthalene	3.843	3.733	2.9	98	0.00
10		Hexachlorobutadiene	0.184	0.181	1.6	99	0.00
11	SURR	2-Methylnaphthalene-d10	0.663	0.642	3.2	100	0.00
12		2-Methylnaphthalene	0.691	0.661	4.3	99	0.00
13	I	Acenaphthene-d10	1.000	1.000	0.0	100	0.00
14	S	2,4,6-Tribromophenol	0.173	0.154	11.0	102	0.00
15	S	2-Fluorobiphenyl	1.626	1.557	4.2	102	0.00
16		Acenaphthylene	1.849	1.709	7.6	99	0.00
17		Acenaphthene	1.113	1.049	5.8	100	0.00
18		Fluorene	1.485	1.427	3.9	101	0.00
19	I	Phenanthrene-d10	1.000	1.000	0.0	102	0.00
20		4-Bromophenyl-phenylether	0.190	0.186	2.1	103	-0.01
21		Hexachlorobenzene	0.196	0.191	2.6	103	-0.01
22		Pentachlorophenol	0.083	0.062	25.3#	90	0.00
23		Phenanthrene	0.980	0.951	3.0	101	-0.01
24		Anthracene	0.933	0.861	7.7	98	0.00
25	SURR	Fluoranthene-d10	5.479	5.078	7.3	97	0.00
26		Fluoranthene	1.412	1.334	5.5	98	0.00
27	I	Chrysene-d12	1.000	1.000	0.0	98	0.00
28		Pyrene	1.247	1.208	3.1	97	0.00
29	S	Terphenyl-d14	0.988	0.940	4.9	96	0.00
30		Benzo(a)anthracene	1.221	1.186	2.9	97	-0.01
31		Chrysene	1.206	1.158	4.0	96	0.00
32		Bis(2-ethylhexyl)phthalate	2.804	2.521	10.1	88	-0.01
33		Indeno(1,2,3-cd)pyrene	1.314	1.276	2.9	96	0.00
34	I	Perylene-d12	1.000	1.000	0.0	97	-0.01
35		Benzo(b)fluoranthene	1.095	1.065	2.7	98	0.00
36		Benzo(k)fluoranthene	1.148	1.142	0.5	100	0.00
37	C	Benzo(a)pyrene	1.064	1.020	4.1	97	0.00
38		Dibenzo(a,h)anthracene	1.006	0.961	4.5	98	-0.01
39		Benzo(g,h,i)perylene	1.029	0.995	3.3	97	-0.02

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