Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN081319\
Data File : BN007130.D
Acq On : 13 Aug 2019 14:07
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Aug 14 04:48:06 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN081019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 13 12:12:02 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	134	0.00
2		n-Nitrosodimethylamine	0.263	0.218	17.1	142	0.00
3	S	2-Fluorophenol	0.961	1.015	-5.6	147	0.00
4	S	Phenol-d6	1.191	1.295	-8.7	157#	0.00
5		bis(2-Chloroethyl)ether	1.052	1.062	-1.0	141	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	143	0.00
7	S	Nitrobenzene-d5	0.323	0.331	-2.5	160#	0.00
8		Naphthalene	1.122	1.051	6.3	143	0.00
9		Hexachlorobutadiene	0.239	0.228	4.6	143	-0.01
10	SURR	2-Methylnaphthalene-d10	0.746	0.687	7.9	142	0.00
11		2-Methylnaphthalene	0.795	0.738	7.2	143	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	137	0.00
13	S	2,4,6-Tribromophenol	0.215	0.180	16.3	157#	0.00
14	S	2-Fluorobiphenyl	0.582	0.334	42.6#	71	0.00
15		Acenaphthylene	2.091	2.351	-12.4	163#	0.00
16		Acenaphthene	1.339	1.294	3.4	141	-0.01
17		Fluorene	2.004	1.764	12.0	145	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	122	0.00
19		4-Bromophenyl-phenylether	0.247	0.251	-1.6	126	-0.01
20		Hexachlorobenzene	0.238	0.243	-2.1	127	0.00
21		Pentachlorophenol	0.089	0.086	3.4	137	0.00
22		Phenanthrene	1.199	1.161	3.2	123	0.00
23		Anthracene	1.095	1.097	-0.2	132	-0.01
24	SURR	Fluoranthene-d10	1.419	1.270	10.5	115	0.00
25		Fluoranthene	1.533	1.411	8.0	117	0.00
26	I	Chrysene-d12	1.000	1.000	0.0	114	-0.01
27		Pyrene	1.403	1.437	-2.4	121	0.00
28	S	Terphenyl-d14	1.090	1.085	0.5	118	-0.01
29		Benzo(a)anthracene	1.463	1.433	2.1	114	0.00
30		Chrysene	1.421	1.403	1.3	117	-0.01
31		Indeno(1,2,3-cd)pyrene	1.566	1.664	-6.3	128	-0.01
32	I	Perylene-d12	1.000	1.000	0.0	126	0.00
33		Benzo(b)fluoranthene	1.372	1.315	4.2	123	0.00
34		Benzo(k)fluoranthene	1.355	1.259	7.1	117	0.00
35	C	Benzo(a)pyrene	1.244	1.220	1.9	129	-0.01
36		Dibenzo(a,h)anthracene	1.241	1.223	1.5	129	-0.01
37		Benzo(g,h,i)perylene	1.276	1.238	3.0	124	-0.01

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