Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN081319\
Data File : BN007147.D
Acq On : 14 Aug 2019 00:20
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 19 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Aug 14 05:00:06 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	153#	0.00
2		n-Nitrosodimethylamine	0.263	0.227	13.7	169#	0.00
3	S	2-Fluorophenol	0.961	1.002	-4.3	167#	0.00
4	S	Phenol-d6	1.191	1.278	-7.3	177#	0.00
5		bis(2-Chloroethyl)ether	1.052	1.026	2.5	156#	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	164#	0.00
7	S	Nitrobenzene-d5	0.323	0.321	0.6	178#	0.00
8		Naphthalene	1.122	1.048	6.6	164#	0.00
9		Hexachlorobutadiene	0.239	0.227	5.0	164#	-0.01
10	SURR	2-Methylnaphthalene-d10	0.746	0.683	8.4	162#	0.00
11		2-Methylnaphthalene	0.795	0.737	7.3	164#	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	159#	0.00
13	S	2,4,6-Tribromophenol	0.215	0.172	20.0	174#	0.00
14	S	2-Fluorobiphenyl	0.582	0.308	47.1#	76	0.00
15		Acenaphthylene	2.091	2.417	-15.6	195#	0.00
16		Acenaphthene	1.339	1.251	6.6	159#	-0.01
17		Fluorene	2.004	1.717	14.3	164#	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	141	0.00
19		4-Bromophenyl-phenylether	0.247	0.245	0.8	141	-0.01
20		Hexachlorobenzene	0.238	0.241	-1.3	144	0.00
21		Pentachlorophenol	0.089	0.079	11.2	145	0.00
22		Phenanthrene	1.199	1.162	3.1	141	0.00
23		Anthracene	1.095	1.073	2.0	148	-0.01
24	SURR	Fluoranthene-d10	1.419	1.259	11.3	131	0.00
25		Fluoranthene	1.533	1.395	9.0	133	0.00
26	I	Chrysene-d12	1.000	1.000	0.0	127	-0.01
27		Pyrene	1.403	1.467	-4.6	137	0.00
28	S	Terphenyl-d14	1.090	1.110	-1.8	134	-0.01
29		Benzo(a)anthracene	1.463	1.495	-2.2	133	-0.01
30		Chrysene	1.421	1.422	-0.1	132	-0.01
31		Indeno(1,2,3-cd)pyrene	1.566	1.706	-8.9	147	-0.02
32	I	Perylene-d12	1.000	1.000	0.0	144	-0.01
33		Benzo(b)fluoranthene	1.372	1.302	5.1	139	-0.01
34		Benzo(k)fluoranthene	1.355	1.273	6.1	136	-0.01
35	C	Benzo(a)pyrene	1.244	1.210	2.7	146	-0.02
36		Dibenzo(a,h)anthracene	1.241	1.224	1.4	148	-0.02
37		Benzo(g,h,i)perylene	1.276	1.223	4.2	140	-0.02

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