Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN081319\
Data File : BN007149.D
Acq On : 14 Aug 2019 02:44
Operator : HP/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Aug 14 05:00:17 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	136	0.00
2		n-Nitrosodimethylamine	0.263	0.233	11.4	153#	0.00
3	S	2-Fluorophenol	0.961	1.014	-5.5	149	0.00
4	S	Phenol-d6	1.191	1.264	-6.1	155#	0.00
5		bis(2-Chloroethyl)ether	1.052	1.050	0.2	141	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	144	0.00
7	S	Nitrobenzene-d5	0.323	0.322	0.3	157#	0.00
8		Naphthalene	1.122	1.050	6.4	144	0.00
9		Hexachlorobutadiene	0.239	0.227	5.0	144	-0.01
10	SURR	2-Methylnaphthalene-d10	0.746	0.687	7.9	143	0.00
11		2-Methylnaphthalene	0.795	0.740	6.9	145	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	138	0.00
13	S	2,4,6-Tribromophenol	0.215	0.174	19.1	153#	0.00
14	S	2-Fluorobiphenyl	0.582	0.285	51.0#	61	0.00
15		Acenaphthylene	2.091	2.296	-9.8	161#	0.00
16		Acenaphthene	1.339	1.276	4.7	141	-0.01
17		Fluorene	2.004	1.708	14.8	142	-0.01
18	I	Phenanthrene-d10	1.000	1.000	0.0	124	0.00
19		4-Bromophenyl-phenylether	0.247	0.246	0.4	125	-0.01
20		Hexachlorobenzene	0.238	0.242	-1.7	128	0.00
21		Pentachlorophenol	0.089	0.080	10.1	130	0.00
22		Phenanthrene	1.199	1.165	2.8	126	0.00
23		Anthracene	1.095	1.070	2.3	131	-0.01
24	SURR	Fluoranthene-d10	1.419	1.254	11.6	115	0.00
25		Fluoranthene	1.533	1.400	8.7	118	0.00
26	I	Chrysene-d12	1.000	1.000	0.0	116	-0.02
27		Pyrene	1.403	1.424	-1.5	121	-0.01
28	S	Terphenyl-d14	1.090	1.072	1.7	118	0.00
29		Benzo(a)anthracene	1.463	1.424	2.7	115	-0.01
30		Chrysene	1.421	1.421	0.0	120	-0.01
31		Indeno(1,2,3-cd)pyrene	1.566	1.682	-7.4	132	-0.02
32	I	Perylene-d12	1.000	1.000	0.0	129	-0.02
33		Benzo(b)fluoranthene	1.372	1.292	5.8	124	-0.02
34		Benzo(k)fluoranthene	1.355	1.280	5.5	122	-0.02
35	C	Benzo(a)pyrene	1.244	1.208	2.9	130	-0.02
36		Dibenzo(a,h)anthracene	1.241	1.222	1.5	132	-0.03
37		Benzo(g,h,i)perylene	1.276	1.222	4.2	125	-0.02

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