Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081019\
Data File : BN007052.D
Acq On : 10 Aug 2019 16:22
Operator : HP/JU
Sample : SSTDICV0.4
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_N
ClientSampleId : ICVBE081019

Quant Time: Aug 10 21:57:19 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-SIM-BN081019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Aug 10 21:44:50 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	104	0.00
2		n-Nitrosodimethylamine	0.263	0.255	3.0	129	0.00
3	S	2-Fluorophenol	0.961	0.905	5.8	103	0.00
4	S	Phenol-d6	1.191	1.074	9.8	101	0.00
5		bis(2-Chloroethyl)ether	1.052	1.018	3.2	106	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	106	0.00
7	S	Nitrobenzene-d5	0.323	0.293	9.3	105	0.00
8		Naphthalene	1.122	1.045	6.9	106	0.00
9		Hexachlorobutadiene	0.239	0.231	3.3	108	-0.01
10	SURR	2-Methylnaphthalene-d10	0.746	0.677	9.2	104	0.00
11		2-Methylnaphthalene	0.795	0.723	9.1	104	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	104	0.00
13	S	2,4,6-Tribromophenol	0.215	0.150	30.2#	99	0.00
14	S	2-Fluorobiphenyl	0.582	0.610	-4.8	99	0.00
15		Acenaphthylene	2.091	1.925	7.9	102	0.00
16		Acenaphthene	1.339	1.256	6.2	104	0.00
17		Fluorene	2.004	1.653	17.5	103	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	103	0.00
19		4-Bromophenyl-phenylether	0.247	0.243	1.6	103	0.00
20		Hexachlorobenzene	0.238	0.241	-1.3	106	0.00
21		Pentachlorophenol	0.089	0.070	21.3	94	0.00
22		Phenanthrene	1.199	1.175	2.0	105	0.00
23		Anthracene	1.095	1.012	7.6	102	0.00
24	SURR	Fluoranthene-d10	1.419	1.339	5.6	102	0.00
25		Fluoranthene	1.533	1.465	4.4	102	0.00
26	I	Chrysene-d12	1.000	1.000	0.0	105	-0.01
27		Pyrene	1.403	1.327	5.4	103	0.00
28	S	Terphenyl-d14	1.090	1.044	4.2	105	0.00
29		Benzo(a)anthracene	1.463	1.395	4.6	103	0.00
30		Chrysene	1.421	1.408	0.9	109	-0.01
31		Indeno(1,2,3-cd)pyrene	1.566	1.385	11.6	99	-0.01
32	I	Perylene-d12	1.000	1.000	0.0	99	0.00
33		Benzo(b)fluoranthene	1.372	1.361	0.8	100	0.00
34		Benzo(k)fluoranthene	1.355	1.383	-2.1	101	0.00
35	C	Benzo(a)pyrene	1.244	1.180	5.1	98	-0.01
36		Dibenzo(a,h)anthracene	1.241	1.193	3.9	99	-0.01
37		Benzo(g,h,i)perylene	1.276	1.247	2.3	98	0.00

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