Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071320\
Data File : BN011158.D
Acq On : 13 Jul 2020 18:11
Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Jul 13 18:43:44 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN071120.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 13 11:25:24 2020
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	83	0.00
2		1,4-Dioxane	0.572	0.567	0.9	77	0.00
3		n-Nitrosodimethylamine	0.288	0.298	-3.5	91	0.00
4	S	2-Fluorophenol	0.986	0.957	2.9	78	0.00
5	S	Phenol-d6	1.115	1.212	-8.7	88	0.00
6		bis(2-Chloroethyl)ether	1.021	1.012	0.9	89	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	88	0.00
8	S	Nitrobenzene-d5	0.355	0.338	4.8	84	0.00
9		Naphthalene	1.190	1.194	-0.3	89	0.00
10		Hexachlorobutadiene	0.295	0.328	-11.2	95	0.00
11	SURR	2-Methylnaphthalene-d10	0.649	0.620	4.5	81	0.00
12		2-Methylnaphthalene	0.749	0.723	3.5	82	0.00
13	I	Acenaphthene-d10	1.000	1.000	0.0	84	0.00
14	S	2,4,6-Tribromophenol	0.178	0.197	-10.7	90	0.00
15	S	2-Fluorobiphenyl	1.874	1.913	-2.1	79	0.00
16		Acenaphthylene	2.036	1.999	1.8	84	0.00
17		Acenaphthene	1.358	1.360	-0.1	83	0.00
18		Fluorene	1.742	1.909	-9.6	80	0.00
19	I	Phenanthrene-d10	1.000	1.000	0.0	85	0.00
20		4,6-Dinitro-2-methylphenol	0.053	0.068	-28.3#	107	0.00
21		4-Bromophenyl-phenylether	0.272	0.315	-15.8	98	0.00
22		Hexachlorobenzene	0.316	0.364	-15.2	96	0.00
23		Atrazine	0.191	0.216	-13.1	95	0.00
24		Pentachlorophenol	0.094	0.101	-7.4	98	0.00
25		Phenanthrene	1.285	1.292	-0.5	85	0.00
26		Anthracene	1.067	1.042	2.3	83	0.00
27	SURR	Fluoranthene-d10	1.261	1.236	2.0	82	0.00
28		Fluoranthene	1.541	1.488	3.4	78	0.00
29	I	Chrysene-d12	1.000	1.000	0.0	84	0.00
30		Pyrene	1.527	1.548	-1.4	75	0.00
31	S	Terphenyl-d14	1.041	1.091	-4.8	83	0.00
32		Benzo(a)anthracene	1.296	1.251	3.5	83	0.00
33		Chrysene	1.584	1.600	-1.0	85	0.00
34		Bis(2-ethylhexyl)phthalate	0.534	0.594	-11.2	95	0.00
35		Indeno(1,2,3-cd)pyrene	1.656	1.833	-10.7	80	0.00
36	I	Perylene-d12	1.000	1.000	0.0	87	0.00
37		Benzo(b)fluoranthene	1.607	1.442	10.3	86	0.00
38		Benzo(k)fluoranthene	1.816	1.914	-5.4	91	0.00
39	C	Benzo(a)pyrene	1.447	1.391	3.9	83	0.00
40		Dibenzo(a,h)anthracene	1.533	1.356	11.5	75	0.00
41		Benzo(g,h,i)perylene	1.683	1.536	8.7	77	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071320\
Data File : BN011158.D
Acq On : 13 Jul 2020 18:11
Operator : CG/JU
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC0.4

Quant Time: Jul 13 18:43:44 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-SIM-BN071120.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 13 11:25:24 2020
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)

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