Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF081619\
Data File : BF116325.D
Acq On : 16 Aug 2019 18:45
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 19 05:30:54 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 07 11:13:18 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	80	0.00
2		1,4-Dioxane	40.000	40.398	-1.0	83	0.00
3		Pyridine	40.000	36.823	7.9	75	0.00
4		n-Nitrosodimethylamine	40.000	36.347	9.1	74	0.00
5	S	2-Fluorophenol	80.000	87.310	-9.1	87	0.00
6		Aniline	40.000	39.375	1.6	78	0.00
7	S	Phenol-d6	80.000	82.988	-3.7	83	0.00
8		2-Chlorophenol	40.000	43.468	-8.7	86	0.00
9		Benzaldehyde	40.000	36.904	7.7	73	0.00
10	C	Phenol	40.000	41.538	-3.8	83	0.00
11		bis(2-Chloroethyl)ether	40.000	43.180	-7.9	87	0.00
12		1,3-Dichlorobenzene	40.000	41.809	-4.5	83	0.00
13	C	1,4-Dichlorobenzene	40.000	42.484	-6.2	84	0.00
14		1,2-Dichlorobenzene	40.000	42.687	-6.7	83	0.00
15		Benzyl Alcohol	40.000	39.984	0.0	78	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	41.998	-5.0	83	0.00
17		2-Methylphenol	40.000	41.126	-2.8	84	0.00
18		Hexachloroethane	40.000	42.674	-6.7	84	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	41.820	-4.6	85	0.00
20		3+4-Methylphenols	40.000	44.107	-10.3	86	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	81	0.00
22		Acetophenone	40.000	42.498	-6.2	85	0.00
23	S	Nitrobenzene-d5	80.000	81.370	-1.7	82	0.00
24		Nitrobenzene	40.000	41.750	-4.4	84	0.00
25		Isophorone	40.000	41.615	-4.0	84	0.00
26	C	2-Nitrophenol	40.000	42.958	-7.4	86	0.00
27		2,4-Dimethylphenol	40.000	41.133	-2.8	82	0.00
28		bis(2-Chloroethoxy)methane	40.000	42.866	-7.2	87	0.00
29	C	2,4-Dichlorophenol	40.000	42.926	-7.3	85	0.00
30		1,2,4-Trichlorobenzene	40.000	41.600	-4.0	83	0.00
31		Naphthalene	40.000	43.236	-8.1	85	0.00
32		Benzoic acid	40.000	33.192	17.0	73	0.00
33		4-Chloroaniline	40.000	41.174	-2.9	82	0.00
34	C	Hexachlorobutadiene	40.000	42.509	-6.3	85	0.00
35		Caprolactam	40.000	39.377	1.6	79	0.00
36	C	4-Chloro-3-methylphenol	40.000	42.134	-5.3	85	0.00
37		2-Methylnaphthalene	40.000	42.435	-6.1	84	0.00
38		1-Methylnaphthalene	40.000	42.314	-5.8	84	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	79	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	43.213	-8.0	84	0.00
41	P	Hexachlorocyclopentadiene	40.000	34.828	12.9	69	0.00
42	S	2,4,6-Tribromophenol	80.000	81.632	-2.0	80	0.00
43	C	2,4,6-Trichlorophenol	40.000	38.460	3.8	75	0.00
44		2,4,5-Trichlorophenol	40.000	37.528	6.2	73	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF081619\
Data File : BF116325.D
Acq On : 16 Aug 2019 18:45
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 19 05:30:54 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 07 11:13:18 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	Amount	Calc.	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	80.000	87.034	-8.8	84	0.00
46		1,1'-Biphenyl	40.000	43.622	-9.1	85	0.00
47		2-Chloronaphthalene	40.000	42.156	-5.4	83	0.00
48		2-Nitroaniline	40.000	41.220	-3.0	81	0.00
49		Acenaphthylene	40.000	43.015	-7.5	83	0.00
50		Dimethylphthalate	40.000	41.383	-3.5	81	0.00
51		2,6-Dinitrotoluene	40.000	41.777	-4.4	82	0.00
52	C	Acenaphthene	40.000	42.643	-6.6	83	0.00
53		3-Nitroaniline	40.000	39.277	1.8	77	0.00
54	P	2,4-Dinitrophenol	40.000	32.810	18.0	64	0.00
55		Dibenzofuran	40.000	41.490	-3.7	80	0.00
56	P	4-Nitrophenol	40.000	37.823	5.4	74	0.00
57		2,4-Dinitrotoluene	40.000	39.429	1.4	76	0.00
58		Fluorene	40.000	44.329	-10.8	86	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	37.720	5.7	74	0.00
60		Diethylphthalate	40.000	42.519	-6.3	82	0.00
61		4-Chlorophenyl-phenylether	40.000	44.649	-11.6	86	0.00
62		4-Nitroaniline	40.000	34.803	13.0	68	0.00
63		Azobenzene	40.000	43.203	-8.0	84	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	76	0.00
65		4,6-Dinitro-2-methylphenol	40.000	36.988	7.5	68	0.00
66	c	n-Nitrosodiphenylamine	40.000	43.801	-9.5	83	0.00
67		4-Bromophenyl-phenylether	40.000	44.252	-10.6	83	0.00
68		Hexachlorobenzene	40.000	43.129	-7.8	81	0.00
69		Atrazine	40.000	40.619	-1.5	77	0.00
70	C	Pentachlorophenol	40.000	35.692	10.8	66	0.00
71		Phenanthrene	40.000	43.285	-8.2	81	0.00
72		Anthracene	40.000	43.653	-9.1	82	0.00
73		Carbazole	40.000	43.632	-9.1	81	0.00
74		Di-n-butylphthalate	40.000	45.307	-13.3	83	0.00
75	C	Fluoranthene	40.000	42.252	-5.6	80	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	68	0.00
77		Benzidine	40.000	35.691	10.8	57	0.00
78		Pyrene	40.000	44.983	-12.5	78	0.00
79	S	Terphenyl-d14	80.000	91.862	-14.8	80	0.00
80		Butylbenzylphthalate	40.000	45.358	-13.4	76	0.00
81		Benzo(a)anthracene	40.000	42.537	-6.3	72	0.00
82		3,3'-Dichlorobenzidine	40.000	40.483	-1.2	67	0.00
83		Chrysene	40.000	43.011	-7.5	73	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	47.162	-17.9	79	0.00
85	c	Di-n-octyl phthalate	40.000	45.876	-14.7	75	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	43.364	-8.4	77	0.00
87	I	Perylene-d12	20.000	20.000	0.0	67	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF081619\
Data File : BF116325.D
Acq On : 16 Aug 2019 18:45
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 19 05:30:54 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 07 11:13:18 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	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	40.000	40.630	-1.6	65	0.00
89	Benzo(k)fluoranthene	40.000	43.366	-8.4	75	0.00
90 C	Benzo(a)pyrene	40.000	42.168	-5.4	70	0.00
91	Dibenzo(a,h)anthracene	40.000	44.924	-12.3	78	0.00
92	Benzo(g,h,i)perylene	40.000	44.325	-10.8	79	0.00

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