Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF082519\
Data File : BF116546.D
Acq On : 26 Aug 2019 00:27
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 26 17:00:56 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 23 13:38:31 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	102	0.00
2		1,4-Dioxane	40.000	35.652	10.9	94	0.00
3		Pyridine	40.000	31.393	21.5	88	0.00
4		n-Nitrosodimethylamine	40.000	36.794	8.0	95	0.01
5	S	2-Fluorophenol	80.000	63.905	20.1	86	0.00
6		Aniline	40.000	36.050	9.9	94	0.00
7	S	Phenol-d6	80.000	65.423	18.2	85	0.01
8		2-Chlorophenol	40.000	32.174	19.6	88	0.00
9		Benzaldehyde	40.000	38.582	3.5	97	0.00
10	C	Phenol	40.000	33.742	15.6	91	0.01
11		bis(2-Chloroethyl)ether	40.000	37.710	5.7	100	0.00
12		1,3-Dichlorobenzene	40.000	38.522	3.7	100	0.00
13	C	1,4-Dichlorobenzene	40.000	40.234	-0.6	101	0.00
14		1,2-Dichlorobenzene	40.000	39.597	1.0	98	0.00
15		Benzyl Alcohol	40.000	32.582	18.5	85	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	34.895	12.8	93	0.00
17		2-Methylphenol	40.000	33.512	16.2	86	0.00
18		Hexachloroethane	40.000	26.404	34.0#	68	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	34.057	14.9	94	0.00
20		3+4-Methylphenols	40.000	31.395	21.5	78	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	94	0.00
22		Acetophenone	40.000	38.547	3.6	95	0.00
23	S	Nitrobenzene-d5	80.000	76.175	4.8	92	0.00
24		Nitrobenzene	40.000	36.591	8.5	88	0.00
25		Isophorone	40.000	35.211	12.0	91	0.00
26	C	2-Nitrophenol	40.000	12.066	69.8#	29	0.00
27		2,4-Dimethylphenol	40.000	30.981	22.5	79	0.00
28		bis(2-Chloroethoxy)methane	40.000	37.537	6.2	96	0.00
29	C	2,4-Dichlorophenol	40.000	27.533	31.2#	67	0.00
30		1,2,4-Trichlorobenzene	40.000	40.239	-0.6	94	0.00
31		Naphthalene	40.000	38.190	4.5	91	0.00
32		Benzoic acid	40.000	6.279	84.3#	14	0.02
33		4-Chloroaniline	40.000	36.700	8.2	87	0.00
34	C	Hexachlorobutadiene	40.000	41.699	-4.2	99	0.00
35		Caprolactam	40.000	25.205	37.0#	63	0.00
36	C	4-Chloro-3-methylphenol	40.000	24.504	38.7#	61	0.01
37		2-Methylnaphthalene	40.000	36.026	9.9	89	0.00
38		1-Methylnaphthalene	40.000	36.136	9.7	89	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	83	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	44.827	-12.1	93	0.00
41	P	Hexachlorocyclopentadiene	40.000	5.369	86.6#	11	0.00
42	S	2,4,6-Tribromophenol	80.000	39.509	50.6#	40	0.00
43	C	2,4,6-Trichlorophenol	40.000	24.602	38.5#	51	0.00
44		2,4,5-Trichlorophenol	40.000	22.624	43.4#	47	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF082519\
Data File : BF116546.D
Acq On : 26 Aug 2019 00:27
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 26 17:00:56 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 23 13:38:31 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	86.656	-8.3	91	0.00
46		1,1'-Biphenyl	40.000	41.798	-4.5	88	0.00
47		2-Chloronaphthalene	40.000	39.945	0.1	84	0.00
48		2-Nitroaniline	40.000	36.389	9.0	77	0.00
49		Acenaphthylene	40.000	37.747	5.6	78	0.00
50		Dimethylphthalate	40.000	42.260	-5.6	89	0.00
51		2,6-Dinitrotoluene	40.000	25.393	36.5#	53	0.00
52	C	Acenaphthene	40.000	36.894	7.8	77	0.00
53		3-Nitroaniline	40.000	41.578	-3.9	85	0.00
54	P	2,4-Dinitrophenol	40.000	8.839	77.9#	1	0.00
55		Dibenzofuran	40.000	37.748	5.6	79	0.00
56	P	4-Nitrophenol	40.000	7.448	81.4#	15	0.00
57		2,4-Dinitrotoluene	40.000	21.376	46.6#	45	0.00
58		Fluorene	40.000	35.440	11.4	76	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	17.331	56.7#	35	0.00
60		Diethylphthalate	40.000	41.171	-2.9	86	0.00
61		4-Chlorophenyl-phenylether	40.000	39.038	2.4	84	0.00
62		4-Nitroaniline	40.000	36.331	9.2	73	0.00
63		Azobenzene	40.000	35.292	11.8	73	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	64	0.00
65		4,6-Dinitro-2-methylphenol	40.000	8.107	79.7#	2	0.00
66	c	n-Nitrosodiphenylamine	40.000	45.617	-14.0	74	0.00
67		4-Bromophenyl-phenylether	40.000	46.046	-15.1	76	0.00
68		Hexachlorobenzene	40.000	43.092	-7.7	72	0.00
69		Atrazine	40.000	37.621	5.9	61	0.00
70	C	Pentachlorophenol	40.000	11.259	71.9#	18	0.00
71		Phenanthrene	40.000	39.956	0.1	64	0.00
72		Anthracene	40.000	39.210	2.0	63	0.00
73		Carbazole	40.000	37.569	6.1	60	0.00
74		Di-n-butylphthalate	40.000	45.046	-12.6	75	0.00
75	C	Fluoranthene	40.000	36.685	8.3	58	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	60	0.00
77		Benzidine	40.000	19.353	51.6#	28	0.00
78		Pyrene	40.000	36.611	8.5	56	0.00
79	S	Terphenyl-d14	80.000	79.765	0.3	63	0.00
80		Butylbenzylphthalate	40.000	40.824	-2.1	64	0.00
81		Benzo(a)anthracene	40.000	38.999	2.5	59	0.00
82		3,3'-Dichlorobenzidine	40.000	32.402	19.0	48	0.00
83		Chrysene	40.000	37.726	5.7	57	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	43.490	-8.7	70	0.00
85	c	Di-n-octyl phthalate	40.000	43.258	-8.1	70	-0.02
86		Indeno(1,2,3-cd)pyrene	40.000	32.175	19.6	59	0.00
87	I	Perylene-d12	20.000	20.000	0.0	56	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF082519\
Data File : BF116546.D
Acq On : 26 Aug 2019 00:27
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 26 17:00:56 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF082319.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 23 13:38:31 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	39.988	0.0	55	0.00
89	Benzo(k)fluoranthene	40.000	42.642	-6.6	57	-0.01
90 C	Benzo(a)pyrene	40.000	37.985	5.0	54	-0.01
91	Dibenzo(a,h)anthracene	40.000	38.184	4.5	59	0.00
92	Benzo(g,h,i)perylene	40.000	38.341	4.1	61	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 5