Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022222\
Data File : BF127008.D
Acq On : 22 Feb 2022 14:53
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Feb 22 23:57:55 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 21 13:16:57 2022
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	76	0.00
2		1,4-Dioxane	40.000	40.473	-1.2	79	-0.01
3		Pyridine	40.000	39.864	0.3	77	-0.02
4		n-Nitrosodimethylamine	40.000	39.908	0.2	75	-0.01
5	S	2-Fluorophenol	80.000	79.570	0.5	77	0.00
6		Aniline	40.000	39.138	2.2	74	0.00
7	S	Phenol-d6	80.000	79.497	0.6	76	0.00
8		2-Chlorophenol	40.000	40.173	-0.4	77	0.00
9		Benzaldehyde	40.000	34.920	12.7	75	0.00
10	C	Phenol	40.000	39.718	0.7	77	0.00
11		bis(2-Chloroethyl)ether	40.000	38.033	4.9	73	0.00
12		1,3-Dichlorobenzene	40.000	40.219	-0.5	79	0.00
13	C	1,4-Dichlorobenzene	40.000	40.203	-0.5	78	0.00
14		1,2-Dichlorobenzene	40.000	40.457	-1.1	78	0.00
15		Benzyl Alcohol	40.000	39.919	0.2	75	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	35.461	11.3	68	0.00
17		2-Methylphenol	40.000	39.620	1.0	76	0.00
18		Hexachloroethane	40.000	40.261	-0.7	78	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.026	2.4	75	0.00
20		3+4-Methylphenols	40.000	40.216	-0.5	76	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	74	0.00
22		Acetophenone	40.000	39.315	1.7	77	0.00
23	S	Nitrobenzene-d5	80.000	79.065	1.2	77	0.00
24		Nitrobenzene	40.000	39.215	2.0	76	0.00
25		Isophorone	40.000	38.121	4.7	73	0.00
26	C	2-Nitrophenol	40.000	40.808	-2.0	77	0.00
27		2,4-Dimethylphenol	40.000	38.971	2.6	76	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.073	4.8	74	0.00
29	C	2,4-Dichlorophenol	40.000	40.270	-0.7	77	0.00
30		1,2,4-Trichlorobenzene	40.000	39.973	0.1	78	0.00
31		Naphthalene	40.000	39.155	2.1	77	0.00
32		Benzoic acid	40.000	41.630	-4.1	74	0.00
33		4-Chloroaniline	40.000	39.852	0.4	77	0.00
34	C	Hexachlorobutadiene	40.000	40.587	-1.5	79	0.00
35		Caprolactam	40.000	39.865	0.3	75	0.00
36	C	4-Chloro-3-methylphenol	40.000	41.006	-2.5	77	0.00
37		2-Methylnaphthalene	40.000	39.532	1.2	76	0.00
38		1-Methylnaphthalene	40.000	39.916	0.2	78	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	76	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.363	-0.9	81	0.00
41	P	Hexachlorocyclopentadiene	40.000	38.264	4.3	72	0.00
42	S	2,4,6-Tribromophenol	80.000	86.745	-8.4	83	0.00
43	C	2,4,6-Trichlorophenol	40.000	39.733	0.7	78	0.00
44		2,4,5-Trichlorophenol	40.000	40.438	-1.1	79	0.00
45	S	2-Fluorobiphenyl	80.000	79.944	0.1	81	0.00
46		1,1'-Biphenyl	40.000	39.031	2.4	78	0.00
47		2-Chloronaphthalene	40.000	38.929	2.7	77	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022222\
Data File : BF127008.D
Acq On : 22 Feb 2022 14:53
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Feb 22 23:57:55 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 21 13:16:57 2022
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)
48		2-Nitroaniline	40.000	40.144	-0.4	79	0.00
49		Acenaphthylene	40.000	39.620	1.0	78	0.00
50		Dimethylphthalate	40.000	39.553	1.1	78	0.00
51		2,6-Dinitrotoluene	40.000	41.119	-2.8	80	0.00
52	C	Acenaphthene	40.000	40.313	-0.8	79	0.00
53		3-Nitroaniline	40.000	39.340	1.6	76	0.00
54	P	2,4-Dinitrophenol	40.000	44.498	-11.2	80	0.00
55		Dibenzofuran	40.000	39.708	0.7	79	0.00
56	P	4-Nitrophenol	40.000	40.958	-2.4	77	0.00
57		2,4-Dinitrotoluene	40.000	41.396	-3.5	79	0.00
58		Fluorene	40.000	40.248	-0.6	80	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	41.868	-4.7	82	0.00
60		Diethylphthalate	40.000	39.892	0.3	79	0.00
61		4-Chlorophenyl-phenylether	40.000	40.583	-1.5	80	0.00
62		4-Nitroaniline	40.000	40.993	-2.5	78	0.00
63		Azobenzene	40.000	38.631	3.4	76	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	79	0.00
65		4,6-Dinitro-2-methylphenol	40.000	44.045	-10.1	82	0.00
66	c	n-Nitrosodiphenylamine	40.000	38.336	4.2	78	0.00
67		4-Bromophenyl-phenylether	40.000	39.374	1.6	80	0.00
68		Hexachlorobenzene	40.000	40.308	-0.8	83	0.00
69		Atrazine	40.000	39.961	0.1	82	0.00
70	C	Pentachlorophenol	40.000	42.718	-6.8	80	0.00
71		Phenanthrene	40.000	39.514	1.2	82	0.00
72		Anthracene	40.000	39.688	0.8	81	0.00
73		Carbazole	40.000	39.890	0.3	82	0.00
74		Di-n-butylphthalate	40.000	39.263	1.8	80	0.00
75	C	Fluoranthene	40.000	41.665	-4.2	86	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	96	0.00
77		Benzidine	40.000	43.964	-9.9	102	0.00
78		Pyrene	40.000	36.567	8.6	87	0.00
79	S	Terphenyl-d14	80.000	75.122	6.1	92	0.00
80		Butylbenzylphthalate	40.000	37.586	6.0	89	0.00
81		Benzo(a)anthracene	40.000	39.996	0.0	97	0.00
82		3,3'-Dichlorobenzidine	40.000	42.793	-7.0	102	0.00
83		Chrysene	40.000	40.129	-0.3	98	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	38.708	3.2	93	0.00
85	c	Di-n-octyl phthalate	40.000	41.873	-4.7	103	0.00
86	I	Perylene-d12	20.000	20.000	0.0	95	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	36.717	8.2	84	0.00
88		Benzo(b)fluoranthene	40.000	41.438	-3.6	102	0.00
89		Benzo(k)fluoranthene	40.000	41.285	-3.2	98	0.00
90	C	Benzo(a)pyrene	40.000	40.612	-1.5	97	0.00
91		Dibenzo(a,h)anthracene	40.000	36.028	9.9	83	0.00
92		Benzo(g,h,i)perylene	40.000	34.733	13.2	79	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022222\
Data File : BF127008.D
Acq On : 22 Feb 2022 14:53
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Feb 22 23:57:55 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF021622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 21 13:16:57 2022
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)

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