Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062222\
Data File : BF128944.D
Acq On : 22 Jun 2022 21:17
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Jun 23 01:33:27 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 22 01:02:29 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	109	-0.01
2		1,4-Dioxane	40.000	41.121	-2.8	106	-0.03
3		Pyridine	40.000	41.507	-3.8	108	-0.02
4		n-Nitrosodimethylamine	40.000	38.799	3.0	101	-0.04
5	S	2-Fluorophenol	80.000	80.656	-0.8	108	0.00
6		Aniline	40.000	41.115	-2.8	113	0.00
7	S	Phenol-d6	80.000	82.553	-3.2	112	0.00
8		2-Chlorophenol	40.000	41.809	-4.5	114	0.00
9		Benzaldehyde	40.000	44.321	-10.8	149	-0.01
10	C	Phenol	40.000	41.406	-3.5	111	0.00
11		bis(2-Chloroethyl)ether	40.000	42.858	-7.1	116	0.00
12		1,3-Dichlorobenzene	40.000	41.698	-4.2	113	-0.01
13	C	1,4-Dichlorobenzene	40.000	41.827	-4.6	115	-0.01
14		1,2-Dichlorobenzene	40.000	42.142	-5.4	117	0.00
15		Benzyl Alcohol	40.000	40.180	-0.4	110	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	43.009	-7.5	121	-0.01
17		2-Methylphenol	40.000	57.635	-44.1#	160	0.00
18		Hexachloroethane	40.000	42.544	-6.4	114	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	43.486	-8.7	118	0.00
20		3+4-Methylphenols	40.000	43.171	-7.9	117	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	115	0.00
22		Acetophenone	40.000	41.570	-3.9	115	0.00
23	S	Nitrobenzene-d5	80.000	84.343	-5.4	118	0.00
24		Nitrobenzene	40.000	42.436	-6.1	118	0.00
25		Isophorone	40.000	42.573	-6.4	119	0.00
26	C	2-Nitrophenol	40.000	45.703	-14.3	123	0.00
27		2,4-Dimethylphenol	40.000	43.191	-8.0	116	0.00
28		bis(2-Chloroethoxy)methane	40.000	42.603	-6.5	117	0.00
29	C	2,4-Dichlorophenol	40.000	42.648	-6.6	116	0.00
30		1,2,4-Trichlorobenzene	40.000	41.483	-3.7	114	0.00
31		Naphthalene	40.000	41.432	-3.6	119	0.00
32		Benzoic acid	40.000	25.007	37.5#	67	0.00
33		4-Chloroaniline	40.000	37.303	6.7	108	0.00
34	C	Hexachlorobutadiene	40.000	40.915	-2.3	119	0.00
35		Caprolactam	40.000	41.875	-4.7	121	0.00
36	C	4-Chloro-3-methylphenol	40.000	41.920	-4.8	121	0.00
37		2-Methylnaphthalene	40.000	41.856	-4.6	121	0.00
38		1-Methylnaphthalene	40.000	41.957	-4.9	124	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	117	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	42.306	-5.8	119	0.00
41	P	Hexachlorocyclopentadiene	40.000	37.618	6.0	102	0.00
42	S	2,4,6-Tribromophenol	80.000	74.222	7.2	101	0.00
43	C	2,4,6-Trichlorophenol	40.000	37.666	5.8	107	0.00
44		2,4,5-Trichlorophenol	40.000	37.726	5.7	106	0.00
45	S	2-Fluorobiphenyl	80.000	82.215	-2.8	117	0.00
46		1,1'-Biphenyl	40.000	41.526	-3.8	119	0.00
47		2-Chloronaphthalene	40.000	42.217	-5.5	122	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062222\
Data File : BF128944.D
Acq On : 22 Jun 2022 21:17
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Jun 23 01:33:27 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 22 01:02:29 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	44.354	-10.9	125	0.00
49		Acenaphthylene	40.000	40.618	-1.5	116	0.00
50		Dimethylphthalate	40.000	42.435	-6.1	120	0.00
51		2,6-Dinitrotoluene	40.000	44.226	-10.6	122	0.00
52	C	Acenaphthene	40.000	38.379	4.1	107	0.00
53		3-Nitroaniline	40.000	43.443	-8.6	120	0.00
54	P	2,4-Dinitrophenol	40.000	35.751	10.6	97	0.00
55		Dibenzofuran	40.000	40.477	-1.2	116	-0.01
56	P	4-Nitrophenol	40.000	34.212	14.5	99	0.00
57		2,4-Dinitrotoluene	40.000	42.689	-6.7	116	0.00
58		Fluorene	40.000	41.602	-4.0	116	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	36.152	9.6	100	0.00
60		Diethylphthalate	40.000	41.784	-4.5	118	0.00
61		4-Chlorophenyl-phenylether	40.000	41.324	-3.3	116	0.00
62		4-Nitroaniline	40.000	39.552	1.1	110	0.00
63		Azobenzene	40.000	42.579	-6.4	120	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	112	0.00
65		4,6-Dinitro-2-methylphenol	40.000	42.215	-5.5	112	0.00
66	c	n-Nitrosodiphenylamine	40.000	42.851	-7.1	116	0.00
67		4-Bromophenyl-phenylether	40.000	43.334	-8.3	117	-0.01
68		Hexachlorobenzene	40.000	42.924	-7.3	115	0.00
69		Atrazine	40.000	37.028	7.4	99	0.00
70	C	Pentachlorophenol	40.000	33.578	16.1	89	0.00
71		Phenanthrene	40.000	41.424	-3.6	114	-0.01
72		Anthracene	40.000	41.284	-3.2	114	0.00
73		Carbazole	40.000	41.403	-3.5	114	0.00
74		Di-n-butylphthalate	40.000	42.299	-5.7	114	0.00
75	C	Fluoranthene	40.000	40.584	-1.5	108	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	91	0.00
77		Benzidine	40.000	34.074	14.8	83	0.00
78		Pyrene	40.000	45.153	-12.9	106	0.00
79	S	Terphenyl-d14	80.000	87.688	-9.6	98	0.00
80		Butylbenzylphthalate	40.000	45.658	-14.1	100	0.00
81		Benzo(a)anthracene	40.000	42.694	-6.7	96	0.00
82		3,3'-Dichlorobenzidine	40.000	47.937	-19.8	109	0.00
83		Chrysene	40.000	42.196	-5.5	96	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	45.425	-13.6	101	0.00
85	c	Di-n-octyl phthalate	40.000	43.881	-9.7	96	0.00
86	I	Perylene-d12	20.000	20.000	0.0	98	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	44.249	-10.6	120	0.00
88		Benzo(b)fluoranthene	40.000	38.732	3.2	92	0.00
89		Benzo(k)fluoranthene	40.000	43.437	-8.6	103	0.00
90	C	Benzo(a)pyrene	40.000	41.548	-3.9	100	0.00
91		Dibenzo(a,h)anthracene	40.000	44.549	-11.4	117	0.00
92		Benzo(g,h,i)perylene	40.000	44.677	-11.7	121	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF062222\
Data File : BF128944.D
Acq On : 22 Jun 2022 21:17
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 21 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Jun 23 01:33:27 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 22 01:02:29 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