Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080922\
Data File : BF129658.D
Acq On : 09 Aug 2022 10:04
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Aug 09 12:26:31 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 05 16:27:34 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	82	0.00
2		1,4-Dioxane	0.554	0.503	9.2	74	-0.02
3		Pyridine	1.538	1.431	7.0	75	-0.02
4		n-Nitrosodimethylamine	0.676	0.643	4.9	77	-0.02
5	S	2-Fluorophenol	1.228	1.168	4.9	80	0.00
6		Aniline	1.918	1.742	9.2	75	0.00
7	S	Phenol-d6	1.543	1.523	1.3	83	0.00
8		2-Chlorophenol	1.341	1.365	-1.8	84	0.00
9		Benzaldehyde	1.048	0.930	11.3	78	0.00
10	C	Phenol	1.643	1.647	-0.2	84	0.00
11		bis(2-Chloroethyl)ether	1.303	1.331	-2.1	85	0.00
12		1,3-Dichlorobenzene	1.439	1.448	-0.6	84	0.00
13	C	1,4-Dichlorobenzene	1.463	1.474	-0.8	84	0.00
14		1,2-Dichlorobenzene	1.345	1.370	-1.9	84	0.00
15		Benzyl Alcohol	1.119	1.192	-6.5	87	0.00
16		2,2'-oxybis(1-Chloropropane	1.959	1.743	11.0	73	0.00
17		2-Methylphenol	1.113	1.102	1.0	83	0.00
18		Hexachloroethane	0.551	0.571	-3.6	86	-0.01
19	P	n-Nitroso-di-n-propylamine	0.934	0.985	-5.5	89	0.00
20		3+4-Methylphenols	1.415	1.417	-0.1	87	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	85	0.00
22		Acetophenone	0.466	0.476	-2.1	91	0.00
23	S	Nitrobenzene-d5	0.366	0.374	-2.2	88	0.00
24		Nitrobenzene	0.377	0.376	0.3	85	0.00
25		Isophorone	0.657	0.673	-2.4	90	0.00
26	C	2-Nitrophenol	0.186	0.194	-4.3	88	-0.01
27		2,4-Dimethylphenol	0.279	0.284	-1.8	88	0.00
28		bis(2-Chloroethoxy)methane	0.381	0.387	-1.6	89	0.00
29	C	2,4-Dichlorophenol	0.288	0.291	-1.0	86	0.00
30		1,2,4-Trichlorobenzene	0.298	0.301	-1.0	87	0.00
31		Naphthalene	1.016	1.014	0.2	87	0.00
32		Benzoic acid	0.202	0.146	27.7#	60	0.00
33		4-Chloroaniline	0.417	0.422	-1.2	89	0.00
34	C	Hexachlorobutadiene	0.170	0.183	-7.6	94	-0.01
35		Caprolactam	0.094	0.095	-1.1	88	0.00
36	C	4-Chloro-3-methylphenol	0.306	0.322	-5.2	91	0.00
37		2-Methylnaphthalene	0.660	0.680	-3.0	90	-0.01
38		1-Methylnaphthalene	0.637	0.662	-3.9	92	-0.01
39	I	Acenaphthene-d10	1.000	1.000	0.0	89	0.00
40		1,2,4,5-Tetrachlorobenzene	0.525	0.544	-3.6	95	0.00
41	P	Hexachlorocyclopentadiene	0.234	0.210	10.3	74	-0.01
42	S	2,4,6-Tribromophenol	0.192	0.202	-5.2	95	0.00
43	C	2,4,6-Trichlorophenol	0.381	0.375	1.6	88	0.00
44		2,4,5-Trichlorophenol	0.415	0.409	1.4	89	0.00
45	S	2-Fluorobiphenyl	1.293	1.250	3.3	98	0.00
46		1,1'-Biphenyl	1.460	1.497	-2.5	93	-0.01
47		2-Chloronaphthalene	1.180	1.175	0.4	91	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080922\
Data File : BF129658.D
Acq On : 09 Aug 2022 10:04
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Aug 09 12:26:31 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 05 16:27:34 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.392	0.390	0.5	88	0.00
49		Acenaphthylene	1.793	1.740	3.0	89	0.00
50		Dimethylphthalate	1.381	1.452	-5.1	96	0.00
51		2,6-Dinitrotoluene	0.309	0.316	-2.3	91	0.00
52	C	Acenaphthene	1.171	1.174	-0.3	90	0.00
53		3-Nitroaniline	0.365	0.351	3.8	85	0.00
54	P	2,4-Dinitrophenol	0.158	0.148	6.3	79	0.00
55		Dibenzofuran	1.614	1.575	2.4	89	0.00
56	P	4-Nitrophenol	0.269	0.232	13.8	76	0.00
57		2,4-Dinitrotoluene	0.404	0.394	2.5	85	0.00
58		Fluorene	1.292	1.312	-1.5	95	0.00
59		2,3,4,6-Tetrachlorophenol	0.321	0.310	3.4	85	0.00
60		Diethylphthalate	1.335	1.404	-5.2	96	0.00
61		4-Chlorophenyl-phenylether	0.569	0.606	-6.5	98	-0.01
62		4-Nitroaniline	0.362	0.338	6.6	84	0.00
63		Azobenzene	1.361	1.355	0.4	90	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	86	0.00
65		4,6-Dinitro-2-methylphenol	0.126	0.131	-4.0	85	0.00
66	c	n-Nitrosodiphenylamine	0.666	0.689	-3.5	92	0.00
67		4-Bromophenyl-phenylether	0.219	0.238	-8.7	96	0.00
68		Hexachlorobenzene	0.237	0.263	-11.0	97	0.00
69		Atrazine	0.202	0.212	-5.0	92	0.00
70	C	Pentachlorophenol	0.129	0.121	6.2	77	0.00
71		Phenanthrene	1.092	1.102	-0.9	90	0.00
72		Anthracene	1.073	1.085	-1.1	91	0.00
73		Carbazole	1.010	0.966	4.4	85	0.00
74		Di-n-butylphthalate	1.203	1.318	-9.6	98	0.00
75	C	Fluoranthene	1.106	1.086	1.8	87	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	79	0.00
77		Benzidine	0.707	0.662	6.4	71	0.00
78		Pyrene	1.672	1.801	-7.7	85	0.00
79	S	Terphenyl-d14	1.060	1.208	-14.0	93	0.00
80		Butylbenzylphthalate	0.725	0.828	-14.2	88	-0.01
81		Benzo(a)anthracene	1.366	1.380	-1.0	81	0.00
82		3,3'-Dichlorobenzidine	0.451	0.441	2.2	77	0.00
83		Chrysene	1.288	1.282	0.5	81	0.00
84		Bis(2-ethylhexyl)phthalate	0.955	1.089	-14.0	91	0.00
85	c	Di-n-octyl phthalate	1.374	1.620	-17.9	95	0.00
86	I	Perylene-d12	1.000	1.000	0.0	86	0.00
87		Indeno(1,2,3-cd)pyrene	1.347	1.460	-8.4	86	0.00
88		Benzo(b)fluoranthene	1.327	1.329	-0.2	89	0.00
89		Benzo(k)fluoranthene	1.300	1.264	2.8	82	0.00
90	C	Benzo(a)pyrene	1.077	1.074	0.3	86	0.00
91		Dibenzo(a,h)anthracene	1.096	1.208	-10.2	87	0.00
92		Benzo(g,h,i)perylene	1.120	1.212	-8.2	84	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080922\
Data File : BF129658.D
Acq On : 09 Aug 2022 10:04
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Aug 09 12:26:31 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF072522.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 05 16:27:34 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	AvgRF	CCRF	%Dev Area%	Dev(min)

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