Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110122\
Data File : BF130902.D
Acq On : 01 Nov 2022 16:01
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Nov 02 08:21:06 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 02 08:18:35 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	88	0.03
2		1,4-Dioxane	0.517	0.458	11.4	74	0.06
3		Pyridine	1.449	1.307	9.8	76	0.06
4		n-Nitrosodimethylamine	0.813	0.773	4.9	81	0.06
5	S	2-Fluorophenol	1.154	1.041	9.8	79	0.03
6		Aniline	1.854	1.653	10.8	76	0.03
7	S	Phenol-d6	1.499	1.319	12.0	78	0.03
8		2-Chlorophenol	1.285	1.210	5.8	81	0.04
9		Benzaldehyde	0.960	0.896	6.7	90	0.03
10	C	Phenol	1.613	1.460	9.5	78	0.02
11		bis(2-Chloroethyl)ether	1.258	1.150	8.6	79	0.03
12		1,3-Dichlorobenzene	1.456	1.352	7.1	80	0.03
13	C	1,4-Dichlorobenzene	1.478	1.366	7.6	80	0.03
14		1,2-Dichlorobenzene	1.369	1.278	6.6	80	0.04
15		Benzyl Alcohol	1.275	1.145	10.2	77	0.03
16		2,2'-oxybis(1-Chloropropane	2.242	2.046	8.7	78	0.02
17		2-Methylphenol	1.098	1.002	8.7	78	0.03
18		Hexachloroethane	0.534	0.537	-0.6	86	0.04
19	P	n-Nitroso-di-n-propylamine	1.020	0.934	8.4	81	0.02
20		3+4-Methylphenols	1.428	1.293	9.5	77	0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	90	0.03
22		Acetophenone	0.485	0.452	6.8	82	0.03
23	S	Nitrobenzene-d5	0.332	0.377	-13.6	92	0.03
24		Nitrobenzene	0.363	0.379	-4.4	87	0.03
25		Isophorone	0.684	0.637	6.9	81	0.03
26	C	2-Nitrophenol	0.123	0.169	-37.4#	110	0.03
27		2,4-Dimethylphenol	0.285	0.264	7.4	79	0.02
28		bis(2-Chloroethoxy)methane	0.382	0.354	7.3	81	0.03
29	C	2,4-Dichlorophenol	0.293	0.278	5.1	81	0.03
30		1,2,4-Trichlorobenzene	0.315	0.304	3.5	84	0.04
31		Naphthalene	1.043	0.958	8.1	80	0.03
32		Benzoic acid	0.177	0.160	9.6	77	0.02
33		4-Chloroaniline	0.412	0.383	7.0	80	0.03
34	C	Hexachlorobutadiene	0.209	0.208	0.5	87	0.03
35		Caprolactam	0.094	0.082	12.8	74	0.03
36	C	4-Chloro-3-methylphenol	0.318	0.306	3.8	82	0.03
37		2-Methylnaphthalene	0.656	0.613	6.6	82	0.03
38		1-Methylnaphthalene	0.639	0.599	6.3	82	0.04
39	I	Acenaphthene-d10	1.000	1.000	0.0	89	0.04
40		1,2,4,5-Tetrachlorobenzene	0.567	0.570	-0.5	88	0.03
41	P	Hexachlorocyclopentadiene	0.298	0.307	-3.0	85	0.04
42	S	2,4,6-Tribromophenol	0.168	0.187	-11.3	90	0.03
43	C	2,4,6-Trichlorophenol	0.394	0.393	0.3	83	0.03
44		2,4,5-Trichlorophenol	0.421	0.427	-1.4	85	0.03
45	S	2-Fluorobiphenyl	1.323	1.249	5.6	87	0.03
46		1,1'-Biphenyl	1.425	1.356	4.8	83	0.04
47		2-Chloronaphthalene	1.143	1.093	4.4	83	0.03

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110122\
Data File : BF130902.D
Acq On : 01 Nov 2022 16:01
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Nov 02 08:21:06 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 02 08:18:35 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	0.315	0.403	-27.9#	99	0.03
49		Acenaphthylene	1.806	1.688	6.5	80	0.03
50		Dimethylphthalate	1.374	1.285	6.5	81	0.03
51		2,6-Dinitrotoluene	0.232	0.283	-22.0	95	0.04
52	C	Acenaphthene	1.116	1.104	1.1	85	0.04
53		3-Nitroaniline	0.268	0.299	-11.6	88	0.03
54	P	2,4-Dinitrophenol	0.092	0.139	-51.1#	132	0.03
55		Dibenzofuran	1.618	1.492	7.8	80	0.04
56	P	4-Nitrophenol	0.212	0.222	-4.7	81	0.02
57		2,4-Dinitrotoluene	0.282	0.372	-31.9#	100	0.03
58		Fluorene	1.274	1.201	5.7	82	0.03
59		2,3,4,6-Tetrachlorophenol	0.347	0.351	-1.2	85	0.03
60		Diethylphthalate	1.351	1.310	3.0	83	0.04
61		4-Chlorophenyl-phenylether	0.610	0.590	3.3	85	0.04
62		4-Nitroaniline	0.271	0.292	-7.7	84	0.04
63		Azobenzene	1.410	1.333	5.5	80	0.03
64	I	Phenanthrene-d10	1.000	1.000	0.0	85	0.03
65		4,6-Dinitro-2-methylphenol	0.073	0.112	-53.4#	128	0.02
66	c	n-Nitrosodiphenylamine	0.624	0.597	4.3	80	0.03
67		4-Bromophenyl-phenylether	0.201	0.210	-4.5	86	0.03
68		Hexachlorobenzene	0.222	0.222	0.0	84	0.03
69		Atrazine	0.180	0.183	-1.7	82	0.03
70	C	Pentachlorophenol	0.126	0.131	-4.0	81	0.03
71		Phenanthrene	1.032	0.970	6.0	79	0.04
72		Anthracene	1.018	0.946	7.1	78	0.04
73		Carbazole	0.912	0.822	9.9	76	0.04
74		Di-n-butylphthalate	1.107	1.102	0.5	83	0.03
75	C	Fluoranthene	1.097	0.979	10.8	75	0.03
76	I	Chrysene-d12	1.000	1.000	0.0	80	0.04
77		Benzidine	0.485	0.389	19.8	62	0.04
78		Pyrene	1.687	1.670	1.0	75	0.04
79	S	Terphenyl-d14	1.091	1.157	-6.0	82	0.04
80		Butylbenzylphthalate	0.607	0.659	-8.6	80	0.04
81		Benzo(a)anthracene	1.326	1.262	4.8	74	0.04
82		3,3'-Dichlorobenzidine	0.358	0.343	4.2	74	0.03
83		Chrysene	1.279	1.172	8.4	72	0.03
84		Bis(2-ethylhexyl)phthalate	0.730	0.801	-9.7	83	0.04
85	c	Di-n-octyl phthalate	1.007	1.127	-11.9	85	0.04
86	I	Perylene-d12	1.000	1.000	0.0	80	0.05
87		Indeno(1,2,3-cd)pyrene	1.328	1.383	-4.1	76	0.07
88		Benzo(b)fluoranthene	1.293	1.207	6.7	75	0.04
89		Benzo(k)fluoranthene	1.260	1.190	5.6	73	0.04
90	C	Benzo(a)pyrene	1.039	0.985	5.2	72	0.05
91		Dibenzo(a,h)anthracene	1.083	1.120	-3.4	74	0.08
92		Benzo(g,h,i)perylene	1.105	1.137	-2.9	73	0.08

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF110122\
Data File : BF130902.D
Acq On : 01 Nov 2022 16:01
Operator : CG\JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Nov 02 08:21:06 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 02 08:18:35 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 = 1