Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092324\
Data File : BP021975.D
Acq On : 23 Sep 2024 09:39
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Sep 23 11:03:53 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 20 01:22:00 2024
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	95	0.00
2		1,4-Dioxane	0.442	0.410	7.2	90	0.00
3		Pyridine	1.179	1.115	5.4	91	0.00
4		n-Nitrosodimethylamine	0.622	0.639	-2.7	100	0.00
5	S	2-Fluorophenol	1.064	1.031	3.1	92	0.00
6		Aniline	1.214	1.215	-0.1	95	0.00
7	S	Phenol-d6	1.382	1.375	0.5	96	0.00
8		2-Chlorophenol	1.138	1.150	-1.1	96	0.00
9		Benzaldehyde	0.722	0.721	0.1	93	0.00
10	C	Phenol	1.372	1.367	0.4	96	0.00
11		bis(2-Chloroethyl)ether	1.034	1.059	-2.4	99	0.00
12		1,3-Dichlorobenzene	1.437	1.456	-1.3	97	0.00
13	C	1,4-Dichlorobenzene	1.457	1.461	-0.3	96	0.00
14		1,2-Dichlorobenzene	1.398	1.391	0.5	95	0.00
15		Benzyl Alcohol	1.060	1.140	-7.5	102	0.00
16		2,2'-oxybis(1-Chloropropane	1.102	1.148	-4.2	100	-0.01
17		2-Methylphenol	0.927	0.982	-5.9	98	0.00
18		Hexachloroethane	0.536	0.553	-3.2	98	-0.01
19	P	n-Nitroso-di-n-propylamine	0.859	0.951	-10.7	103	0.00
20		3+4-Methylphenols	1.309	1.371	-4.7	98	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	100	0.00
22		Acetophenone	0.488	0.464	4.9	100	-0.01
23	S	Nitrobenzene-d5	0.378	0.375	0.8	102	0.00
24		Nitrobenzene	0.384	0.380	1.0	101	0.00
25		Isophorone	0.623	0.623	0.0	101	-0.01
26	C	2-Nitrophenol	0.168	0.166	1.2	99	0.00
27		2,4-Dimethylphenol	0.194	0.197	-1.5	102	0.00
28		bis(2-Chloroethoxy)methane	0.374	0.368	1.6	100	0.00
29	C	2,4-Dichlorophenol	0.330	0.335	-1.5	100	0.00
30		1,2,4-Trichlorobenzene	0.411	0.406	1.2	101	-0.01
31		Naphthalene	0.998	0.990	0.8	102	0.00
32		Benzoic acid	0.227	0.234	-3.1	100	0.00
33		4-Chloroaniline	0.360	0.367	-1.9	102	0.00
34	C	Hexachlorobutadiene	0.321	0.327	-1.9	106	0.00
35		Caprolactam	0.098	0.099	-1.0	100	0.00
36	C	4-Chloro-3-methylphenol	0.354	0.373	-5.4	106	0.00
37		2-Methylnaphthalene	0.730	0.748	-2.5	105	-0.01
38		1-Methylnaphthalene	0.721	0.753	-4.4	107	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	112	0.00
40		1,2,4,5-Tetrachlorobenzene	0.700	0.673	3.9	109	0.00
41	P	Hexachlorocyclopentadiene	0.260	0.251	3.5	110	0.00
42	S	2,4,6-Tribromophenol	0.371	0.410	-10.5	127	-0.02
43	C	2,4,6-Trichlorophenol	0.429	0.415	3.3	107	-0.01
44		2,4,5-Trichlorophenol	0.482	0.474	1.7	111	0.00
45	S	2-Fluorobiphenyl	1.370	1.339	2.3	113	0.00
46		1,1'-Biphenyl	1.364	1.330	2.5	112	-0.01
47		2-Chloronaphthalene	1.103	1.070	3.0	111	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092324\
Data File : BP021975.D
Acq On : 23 Sep 2024 09:39
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Sep 23 11:03:53 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 20 01:22:00 2024
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	0.301	0.315	-4.7	111	0.00
49		Acenaphthylene	1.556	1.567	-0.7	112	0.00
50		Dimethylphthalate	1.451	1.438	0.9	112	-0.02
51		2,6-Dinitrotoluene	0.323	0.322	0.3	111	0.00
52	C	Acenaphthene	1.013	1.021	-0.8	115	0.00
53		3-Nitroaniline	0.288	0.298	-3.5	111	0.00
54	P	2,4-Dinitrophenol	0.189	0.197	-4.2	113	0.00
55		Dibenzofuran	1.702	1.761	-3.5	118	0.00
56	P	4-Nitrophenol	0.251	0.261	-4.0	109	0.00
57		2,4-Dinitrotoluene	0.457	0.486	-6.3	116	0.00
58		Fluorene	1.421	1.482	-4.3	119	0.00
59		2,3,4,6-Tetrachlorophenol	0.465	0.481	-3.4	117	0.00
60		Diethylphthalate	1.474	1.546	-4.9	121	-0.01
61		4-Chlorophenyl-phenylether	0.828	0.873	-5.4	123	0.00
62		4-Nitroaniline	0.317	0.344	-8.5	117	0.00
63		Azobenzene	1.220	1.333	-9.3	124	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	124	-0.02
65		4,6-Dinitro-2-methylphenol	0.116	0.116	0.0	121	-0.02
66	c	n-Nitrosodiphenylamine	0.498	0.485	2.6	120	0.00
67		4-Bromophenyl-phenylether	0.233	0.229	1.7	124	-0.01
68		Hexachlorobenzene	0.291	0.288	1.0	125	0.00
69		Atrazine	0.178	0.148	16.9	114	-0.02
70	C	Pentachlorophenol	0.196	0.197	-0.5	123	-0.01
71		Phenanthrene	0.960	0.929	3.2	121	-0.02
72		Anthracene	0.929	0.933	-0.4	125	0.00
73		Carbazole	0.898	0.906	-0.9	123	-0.01
74		Di-n-butylphthalate	1.054	1.087	-3.1	131	0.00
75	C	Fluoranthene	1.340	1.391	-3.8	128	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	131	-0.01
77		Benzidine	0.244	0.390	-59.8#	127	-0.02
78		Pyrene	1.210	1.167	3.6	126	0.00
79	S	Terphenyl-d14	1.065	1.042	2.2	129	0.00
80		Butylbenzylphthalate	0.432	0.428	0.9	130	-0.01
81		Benzo(a)anthracene	1.285	1.254	2.4	130	-0.01
82		3,3'-Dichlorobenzidine	0.466	0.463	0.6	122	0.00
83		Chrysene	1.210	1.187	1.9	132	-0.02
84		Bis(2-ethylhexyl)phthalate	0.633	0.638	-0.8	138	0.00
85	c	Di-n-octyl phthalate	1.076	1.079	-0.3	136	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	132	-0.04
87		Indeno(1,2,3-cd)pyrene	1.330	1.311	1.4	130	-0.05
88		Benzo(b)fluoranthene	1.183	1.189	-0.5	130	-0.03
89		Benzo(k)fluoranthene	1.119	1.093	2.3	131	-0.02
90	C	Benzo(a)pyrene	1.019	1.001	1.8	129	-0.02
91		Dibenzo(a,h)anthracene	1.098	1.063	3.2	129	-0.04
92		Benzo(g,h,i)perylene	1.125	1.128	-0.3	131	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092324\
Data File : BP021975.D
Acq On : 23 Sep 2024 09:39
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Sep 23 11:03:53 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 20 01:22:00 2024
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