Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021622\
Data File : BP009171.D
Acq On : 16 Feb 2022 10:38
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Feb 16 14:30:53 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 16 14:30:24 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	123	0.00
2		1,4-Dioxane	0.369	0.359	2.7	122	0.00
3		Pyridine	1.079	1.021	5.4	117	0.00
4		n-Nitrosodimethylamine	0.413	0.365	11.6	105	0.00
5	S	2-Fluorophenol	1.121	0.994	11.3	111	0.00
6		Aniline	1.677	1.511	9.9	115	0.00
7	S	Phenol-d6	1.477	1.314	11.0	113	0.00
8		2-Chlorophenol	1.265	1.123	11.2	113	0.00
9		Benzaldehyde	0.748	0.666	11.0	116	0.00
10	C	Phenol	1.485	1.319	11.2	112	0.00
11		bis(2-Chloroethyl)ether	1.153	1.056	8.4	118	0.00
12		1,3-Dichlorobenzene	1.463	1.335	8.7	117	0.00
13	C	1,4-Dichlorobenzene	1.494	1.349	9.7	115	0.00
14		1,2-Dichlorobenzene	1.448	1.280	11.6	113	0.00
15		Benzyl Alcohol	0.940	0.964	-2.6	128	0.00
16		2,2'-oxybis(1-Chloropropane	1.300	1.177	9.5	118	0.00
17		2-Methylphenol	0.996	0.929	6.7	119	0.00
18		Hexachloroethane	0.495	0.463	6.5	121	0.00
19	P	n-Nitroso-di-n-propylamine	0.846	0.804	5.0	123	0.00
20		3+4-Methylphenols	1.363	1.292	5.2	121	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	129	0.00
22		Acetophenone	0.468	0.435	7.1	120	0.00
23	S	Nitrobenzene-d5	0.355	0.328	7.6	119	0.00
24		Nitrobenzene	0.335	0.319	4.8	123	0.00
25		Isophorone	0.589	0.600	-1.9	138	0.00
26	C	2-Nitrophenol	0.173	0.183	-5.8	134	0.00
27		2,4-Dimethylphenol	0.259	0.252	2.7	127	0.00
28		bis(2-Chloroethoxy)methane	0.400	0.377	5.8	126	0.00
29	C	2,4-Dichlorophenol	0.324	0.327	-0.9	131	0.00
30		1,2,4-Trichlorobenzene	0.381	0.371	2.6	128	0.00
31		Naphthalene	1.020	0.963	5.6	124	0.00
32		Benzoic acid	0.202	0.203	-0.5	138	0.00
33		4-Chloroaniline	0.412	0.409	0.7	131	0.00
34	C	Hexachlorobutadiene	0.234	0.234	0.0	132	0.00
35		Caprolactam	0.095	0.106	-11.6	157#	0.00
36	C	4-Chloro-3-methylphenol	0.309	0.315	-1.9	139	0.00
37		2-Methylnaphthalene	0.722	0.694	3.9	130	0.00
38		1-Methylnaphthalene	0.705	0.681	3.4	131	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	148	0.00
40		1,2,4,5-Tetrachlorobenzene	0.637	0.600	5.8	136	0.00
41	P	Hexachlorocyclopentadiene	0.186	0.217	-16.7	168#	0.00
42	S	2,4,6-Tribromophenol	0.267	0.286	-7.1	167#	0.00
43	C	2,4,6-Trichlorophenol	0.422	0.417	1.2	142	0.00
44		2,4,5-Trichlorophenol	0.453	0.444	2.0	142	0.00
45	S	2-Fluorobiphenyl	1.429	1.309	8.4	133	0.00
46		1,1'-Biphenyl	1.469	1.363	7.2	136	0.00
47		2-Chloronaphthalene	1.168	1.069	8.5	134	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021622\
Data File : BP009171.D
Acq On : 16 Feb 2022 10:38
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Feb 16 14:30:53 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 16 14:30:24 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	0.271	0.269	0.7	146	0.00
49		Acenaphthylene	1.760	1.644	6.6	138	0.00
50		Dimethylphthalate	1.426	1.383	3.0	151#	0.00
51		2,6-Dinitrotoluene	0.297	0.297	0.0	151#	0.00
52	C	Acenaphthene	1.071	0.992	7.4	138	0.00
53		3-Nitroaniline	0.304	0.311	-2.3	153#	0.00
54	P	2,4-Dinitrophenol	0.157	0.162	-3.2	158#	0.00
55		Dibenzofuran	1.805	1.690	6.4	142	0.00
56	P	4-Nitrophenol	0.218	0.317	-45.4#	225#	0.00
57		2,4-Dinitrotoluene	0.388	0.414	-6.7	162#	0.00
58		Fluorene	1.386	1.312	5.3	144	0.00
59		2,3,4,6-Tetrachlorophenol	0.396	0.401	-1.3	155#	0.00
60		Diethylphthalate	1.348	1.364	-1.2	161#	0.00
61		4-Chlorophenyl-phenylether	0.753	0.742	1.5	151#	0.00
62		4-Nitroaniline	0.306	0.316	-3.3	154#	0.00
63		Azobenzene	1.166	1.135	2.7	150	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	170#	0.00
65		4,6-Dinitro-2-methylphenol	0.123	0.123	0.0	173#	0.00
66	c	n-Nitrosodiphenylamine	0.607	0.553	8.9	152#	0.00
67		4-Bromophenyl-phenylether	0.232	0.222	4.3	163#	0.00
68		Hexachlorobenzene	0.260	0.251	3.5	165#	0.00
69		Atrazine	0.201	0.208	-3.5	178#	0.00
70	C	Pentachlorophenol	0.137	0.143	-4.4	176#	0.00
71		Phenanthrene	1.083	1.000	7.7	157#	0.00
72		Anthracene	1.060	1.001	5.6	160#	0.00
73		Carbazole	1.016	0.969	4.6	161#	0.00
74		Di-n-butylphthalate	1.029	1.104	-7.3	188#	0.00
75	C	Fluoranthene	1.214	1.216	-0.2	170#	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	178#	0.00
77		Benzidine	0.418	0.358	14.4	128	0.00
78		Pyrene	1.423	1.217	14.5	170#	0.00
79	S	Terphenyl-d14	1.112	0.959	13.8	171#	0.00
80		Butylbenzylphthalate	0.489	0.506	-3.5	202#	0.00
81		Benzo(a)anthracene	1.289	1.220	5.4	169#	0.00
82		3,3'-Dichlorobenzidine	0.447	0.452	-1.1	169#	0.00
83		Chrysene	1.242	1.144	7.9	167#	0.00
84		Bis(2-ethylhexyl)phthalate	0.647	0.718	-11.0	212#	0.00
85	c	Di-n-octyl phthalate	1.035	1.263	-22.0#	209#	0.00
86	I	Perylene-d12	1.000	1.000	0.0	164#	0.00
87		Indeno(1,2,3-cd)pyrene	1.486	1.318	11.3	135	0.00
88		Benzo(b)fluoranthene	1.231	1.145	7.0	158#	0.00
89		Benzo(k)fluoranthene	1.228	1.138	7.3	156#	0.00
90	C	Benzo(a)pyrene	1.027	0.964	6.1	153#	0.00
91		Dibenzo(a,h)anthracene	1.221	1.087	11.0	135	0.00
92		Benzo(g,h,i)perylene	1.216	1.033	15.0	129	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021622\
Data File : BP009171.D
Acq On : 16 Feb 2022 10:38
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Feb 16 14:30:53 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 16 14:30:24 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