Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022624\
Data File : BP019884.D
Acq On : 26 Feb 2024 18:03
Operator : MA/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ICVBP022624

Quant Time: Feb 27 02:06:46 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP022624.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 27 02:05:05 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	89	0.00
2		1,4-Dioxane	0.421	0.413	1.9	87	0.00
3		Pyridine	1.289	1.338	-3.8	86	0.00
4		n-Nitrosodimethylamine	0.531	0.561	-5.6	88	0.00
5	S	2-Fluorophenol	1.218	1.165	4.4	87	0.00
6		Aniline	2.225	2.140	3.8	87	0.00
7	S	Phenol-d6	1.814	1.752	3.4	86	0.00
8		2-Chlorophenol	1.314	1.299	1.1	87	0.00
9		Benzaldehyde	1.035	0.976	5.7	85	0.00
10	C	Phenol	1.809	1.710	5.5	85	0.00
11		bis(2-Chloroethyl)ether	1.425	1.336	6.2	84	0.00
12		1,3-Dichlorobenzene	1.427	1.428	-0.1	88	0.00
13	C	1,4-Dichlorobenzene	1.445	1.466	-1.5	89	0.00
14		1,2-Dichlorobenzene	1.449	1.442	0.5	89	0.00
15		Benzyl Alcohol	1.600	1.539	3.8	85	0.00
16		2,2'-oxybis(1-Chloropropane	1.521	1.320	13.2	81	0.00
17		2-Methylphenol	1.353	1.297	4.1	86	0.00
18		Hexachloroethane	0.560	0.565	-0.9	89	0.00
19	P	n-Nitroso-di-n-propylamine	1.370	1.327	3.1	84	0.00
20		3+4-Methylphenols	1.944	1.835	5.6	86	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	88	0.00
22		Acetophenone	0.621	0.558	10.1	86	0.00
23	S	Nitrobenzene-d5	0.495	0.462	6.7	86	0.00
24		Nitrobenzene	0.480	0.452	5.8	87	0.00
25		Isophorone	0.862	0.829	3.8	85	0.00
26	C	2-Nitrophenol	0.192	0.198	-3.1	89	0.00
27		2,4-Dimethylphenol	0.317	0.319	-0.6	88	0.00
28		bis(2-Chloroethoxy)methane	0.481	0.464	3.5	84	0.00
29	C	2,4-Dichlorophenol	0.353	0.364	-3.1	89	0.00
30		1,2,4-Trichlorobenzene	0.388	0.397	-2.3	89	0.00
31		Naphthalene	1.071	1.065	0.6	87	0.00
32		Benzoic acid	0.268	0.283	-5.6	90	0.00
33		4-Chloroaniline	0.487	0.491	-0.8	87	0.00
34	C	Hexachlorobutadiene	0.282	0.292	-3.5	90	0.00
35		Caprolactam	0.130	0.132	-1.5	88	0.00
36	C	4-Chloro-3-methylphenol	0.436	0.434	0.5	86	0.00
37		2-Methylnaphthalene	0.831	0.837	-0.7	88	0.00
38		1-Methylnaphthalene	0.787	0.785	0.3	88	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	100	0.00
40		1,2,4,5-Tetrachlorobenzene	0.683	0.620	9.2	90	0.00
41	P	Hexachlorocyclopentadiene	0.354	0.338	4.5	90	0.00
42	S	2,4,6-Tribromophenol	0.343	0.337	1.7	94	0.00
43	C	2,4,6-Trichlorophenol	0.451	0.417	7.5	91	0.00
44		2,4,5-Trichlorophenol	0.537	0.500	6.9	92	0.00
45	S	2-Fluorobiphenyl	1.480	1.341	9.4	90	0.00
46		1,1'-Biphenyl	1.432	1.283	10.4	89	0.00
47		2-Chloronaphthalene	1.112	1.000	10.1	89	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022624\
Data File : BP019884.D
Acq On : 26 Feb 2024 18:03
Operator : MA/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : ICVBP022624

Quant Time: Feb 27 02:06:46 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP022624.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 27 02:05:05 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	0.402	0.357	11.2	86	0.00
49		Acenaphthylene	1.801	1.790	0.6	98	0.00
50		Dimethylphthalate	1.633	1.480	9.4	89	0.00
51		2,6-Dinitrotoluene	0.351	0.340	3.1	95	0.00
52	C	Acenaphthene	1.081	1.061	1.9	97	0.00
53		3-Nitroaniline	0.348	0.375	-7.8	105	0.00
54	P	2,4-Dinitrophenol	0.229	0.215	6.1	90	0.00
55		Dibenzofuran	1.845	1.695	8.1	90	0.00
56	P	4-Nitrophenol	0.272	0.259	4.8	88	0.00
57		2,4-Dinitrotoluene	0.500	0.467	6.6	90	0.00
58		Fluorene	1.497	1.392	7.0	91	0.00
59		2,3,4,6-Tetrachlorophenol	0.493	0.473	4.1	92	0.00
60		Diethylphthalate	1.613	1.471	8.8	88	0.00
61		4-Chlorophenyl-phenylether	0.856	0.808	5.6	92	0.00
62		4-Nitroaniline	0.362	0.347	4.1	89	0.00
63		Azobenzene	1.502	1.355	9.8	86	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	91	0.00
65		4,6-Dinitro-2-methylphenol	0.136	0.145	-6.6	91	0.00
66	c	n-Nitrosodiphenylamine	0.574	0.575	-0.2	89	0.00
67		4-Bromophenyl-phenylether	0.247	0.253	-2.4	92	0.00
68		Hexachlorobenzene	0.267	0.277	-3.7	94	0.00
69		Atrazine	0.240	0.243	-1.3	89	0.00
70	C	Pentachlorophenol	0.197	0.206	-4.6	90	0.00
71		Phenanthrene	1.068	1.067	0.1	89	0.00
72		Anthracene	1.090	1.101	-1.0	90	0.00
73		Carbazole	0.965	0.974	-0.9	89	0.00
74		Di-n-butylphthalate	1.207	1.190	1.4	86	0.00
75	C	Fluoranthene	1.385	1.411	-1.9	89	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	89	0.00
77		Benzidine	0.574	0.592	-3.1	83	0.00
78		Pyrene	1.374	1.356	1.3	89	0.00
79	S	Terphenyl-d14	1.269	1.259	0.8	91	0.00
80		Butylbenzylphthalate	0.549	0.530	3.5	86	0.00
81		Benzo(a)anthracene	1.420	1.421	-0.1	89	0.00
82		3,3'-Dichlorobenzidine	0.544	0.555	-2.0	90	0.00
83		Chrysene	1.330	1.346	-1.2	91	0.00
84		Bis(2-ethylhexyl)phthalate	0.789	0.756	4.2	84	0.00
85	c	Di-n-octyl phthalate	1.313	1.271	3.2	84	0.00
86	I	Perylene-d12	1.000	1.000	0.0	90	0.00
87		Indeno(1,2,3-cd)pyrene	1.408	1.448	-2.8	90	0.00
88		Benzo(b)fluoranthene	1.226	1.226	0.0	92	0.00
89		Benzo(k)fluoranthene	1.203	1.228	-2.1	89	0.00
90	C	Benzo(a)pyrene	1.172	1.189	-1.5	90	0.00
91		Dibenzo(a,h)anthracene	1.184	1.214	-2.5	90	0.00
92		Benzo(g,h,i)perylene	1.143	1.169	-2.3	89	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022624\
Data File : BP019884.D
Acq On : 26 Feb 2024 18:03
Operator : MA/JU
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ICVBP022624

Quant Time: Feb 27 02:06:46 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP022624.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 27 02:05:05 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