Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012924\
Data File : BP019470.D
Acq On : 29 Jan 2024 17:38
Operator : MA/JU
Sample : SSTDICV040
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ICVBP012924

Quant Time: Jan 30 01:00:53 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP012924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 30 00:59:08 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	92	0.00
2		1,4-Dioxane	40.000	37.228	6.9	89	0.00
3		Pyridine	40.000	37.682	5.8	92	0.00
4		n-Nitrosodimethylamine	40.000	38.706	3.2	92	0.00
5	S	2-Fluorophenol	80.000	75.478	5.7	91	0.00
6		Aniline	40.000	37.220	7.0	93	0.00
7	S	Phenol-d6	80.000	73.981	7.5	91	0.00
8		2-Chlorophenol	40.000	36.933	7.7	91	0.00
9		Benzaldehyde	40.000	38.988	2.5	91	0.00
10	C	Phenol	40.000	37.728	5.7	93	0.00
11		bis(2-Chloroethyl)ether	40.000	36.852	7.9	92	0.00
12		1,3-Dichlorobenzene	40.000	37.180	7.1	91	0.00
13	C	1,4-Dichlorobenzene	40.000	36.733	8.2	90	0.00
14		1,2-Dichlorobenzene	40.000	37.250	6.9	91	0.00
15		Benzyl Alcohol	40.000	37.000	7.5	93	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	35.821	10.4	89	0.00
17		2-Methylphenol	40.000	37.020	7.4	94	0.00
18		Hexachloroethane	40.000	37.057	7.4	90	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	36.174	9.6	93	0.00
20		3+4-Methylphenols	40.000	37.255	6.9	95	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	94	0.00
22		Acetophenone	40.000	37.771	5.6	94	0.00
23	S	Nitrobenzene-d5	80.000	77.057	3.7	94	0.00
24		Nitrobenzene	40.000	38.120	4.7	94	0.00
25		Isophorone	40.000	36.989	7.5	94	0.00
26	C	2-Nitrophenol	40.000	40.385	-1.0	96	0.00
27		2,4-Dimethylphenol	40.000	37.923	5.2	94	0.00
28		bis(2-Chloroethoxy)methane	40.000	37.224	6.9	93	0.00
29	C	2,4-Dichlorophenol	40.000	38.052	4.9	94	0.00
30		1,2,4-Trichlorobenzene	40.000	37.989	5.0	93	0.00
31		Naphthalene	40.000	37.517	6.2	94	0.00
32		Benzoic acid	40.000	40.713	-1.8	102	0.00
33		4-Chloroaniline	40.000	37.951	5.1	95	0.00
34	C	Hexachlorobutadiene	40.000	37.600	6.0	92	0.00
35		Caprolactam	40.000	39.167	2.1	103	0.00
36	C	4-Chloro-3-methylphenol	40.000	38.492	3.8	98	0.00
37		2-Methylnaphthalene	40.000	37.005	7.5	94	0.00
38		1-Methylnaphthalene	40.000	37.170	7.1	94	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	98	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	37.850	5.4	95	0.00
41	P	Hexachlorocyclopentadiene	40.000	37.540	6.2	91	0.00
42	S	2,4,6-Tribromophenol	80.000	80.976	-1.2	107	0.00
43	C	2,4,6-Trichlorophenol	40.000	39.121	2.2	98	0.00
44		2,4,5-Trichlorophenol	40.000	38.702	3.2	97	0.00
45	S	2-Fluorobiphenyl	80.000	75.441	5.7	96	0.00
46		1,1'-Biphenyl	40.000	38.278	4.3	98	0.00
47		2-Chloronaphthalene	40.000	38.204	4.5	97	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012924\
Data File : BP019470.D
Acq On : 29 Jan 2024 17:38
Operator : MA/JU
Sample : SSTDICV040
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ICVBP012924

Quant Time: Jan 30 01:00:53 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP012924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 30 00:59:08 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	Amount	Calc.	%Dev	Area%	Dev(min)
48		2-Nitroaniline	40.000	40.295	-0.7	101	0.00
49		Acenaphthylene	40.000	38.320	4.2	99	0.00
50		Dimethylphthalate	40.000	38.539	3.7	101	0.00
51		2,6-Dinitrotoluene	40.000	38.885	2.8	101	0.00
52	C	Acenaphthene	40.000	38.221	4.4	98	0.00
53		3-Nitroaniline	40.000	40.379	-0.9	105	0.00
54	P	2,4-Dinitrophenol	40.000	40.219	-0.5	111	0.00
55		Dibenzofuran	40.000	38.256	4.4	99	0.00
56	P	4-Nitrophenol	40.000	42.205	-5.5	109	0.00
57		2,4-Dinitrotoluene	40.000	40.608	-1.5	107	0.00
58		Fluorene	40.000	38.466	3.8	101	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	39.190	2.0	103	0.00
60		Diethylphthalate	40.000	38.774	3.1	104	0.00
61		4-Chlorophenyl-phenylether	40.000	38.095	4.8	100	0.00
62		4-Nitroaniline	40.000	41.019	-2.5	109	0.00
63		Azobenzene	40.000	38.455	3.9	100	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	109	0.00
65		4,6-Dinitro-2-methylphenol	40.000	39.805	0.5	115	0.00
66	c	n-Nitrosodiphenylamine	40.000	37.041	7.4	104	0.00
67		4-Bromophenyl-phenylether	40.000	37.763	5.6	107	0.00
68		Hexachlorobenzene	40.000	37.341	6.6	105	0.00
69		Atrazine	40.000	38.814	3.0	111	0.00
70	C	Pentachlorophenol	40.000	40.599	-1.5	112	0.00
71		Phenanthrene	40.000	37.545	6.1	108	0.00
72		Anthracene	40.000	37.932	5.2	107	0.00
73		Carbazole	40.000	38.833	2.9	113	0.00
74		Di-n-butylphthalate	40.000	38.987	2.5	117	0.00
75	C	Fluoranthene	40.000	39.666	0.8	122	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	150	0.00
77		Benzidine	40.000	43.734	-9.3	176	0.00
78		Pyrene	40.000	35.288	11.8	126	0.00
79	S	Terphenyl-d14	80.000	71.626	10.5	130	0.00
80		Butylbenzylphthalate	40.000	37.567	6.1	148	0.00
81		Benzo(a)anthracene	40.000	37.704	5.7	148	0.00
82		3,3'-Dichlorobenzidine	40.000	38.513	3.7	157	0.00
83		Chrysene	40.000	38.165	4.6	149	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	38.650	3.4	163	0.00
85	c	Di-n-octyl phthalate	40.000	39.208	2.0	170	0.00
86	I	Perylene-d12	20.000	20.000	0.0	143	0.00
87		Indeno(1,2,3-cd)pyrene	40.000	33.475	16.3	112	0.00
88		Benzo(b)fluoranthene	40.000	40.205	-0.5	159	0.00
89		Benzo(k)fluoranthene	40.000	38.344	4.1	148	0.00
90	C	Benzo(a)pyrene	40.000	38.422	3.9	146	0.00
91		Dibenzo(a,h)anthracene	40.000	33.507	16.2	113	0.00
92		Benzo(g,h,i)perylene	40.000	32.724	18.2	105	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP012924\
Data File : BP019470.D
Acq On : 29 Jan 2024 17:38
Operator : MA/JU
Sample : SSTDICV040
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
ICVBP012924

Quant Time: Jan 30 01:00:53 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP012924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 30 00:59:08 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	Amount	Calc.	%Dev	Area%	Dev(min)

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