Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081721\
Data File : BF125073.D
Acq On : 17 Aug 2021 17:30
Operator : JU/CG
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF081721

Quant Time: Aug 17 18:18:39 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 17 16:45:36 2021
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	91	0.00
2		1,4-Dioxane	0.680	0.635	6.6	82	-0.02
3		Pyridine	1.844	1.718	6.8	81	-0.01
4		n-Nitrosodimethylamine	0.989	0.916	7.4	80	-0.02
5	S	2-Fluorophenol	1.339	1.249	6.7	85	0.00
6		Aniline	2.079	1.864	10.3	79	0.00
7	S	Phenol-d6	1.647	1.502	8.8	83	0.00
8		2-Chlorophenol	1.340	1.253	6.5	83	0.00
9		Benzaldehyde	1.246	1.093	12.3	82	0.00
10	C	Phenol	1.841	1.600	13.1	80	0.00
11		bis(2-Chloroethyl)ether	1.402	1.290	8.0	83	0.00
12		1,3-Dichlorobenzene	1.428	1.319	7.6	83	0.00
13	C	1,4-Dichlorobenzene	1.431	1.343	6.1	84	0.00
14		1,2-Dichlorobenzene	1.378	1.227	11.0	84	0.00
15		Benzyl Alcohol	1.343	1.218	9.3	78	0.00
16		2,2'-oxybis(1-Chloropropane	3.129	2.869	8.3	81	0.00
17		2-Methylphenol	1.202	1.099	8.6	81	0.00
18		Hexachloroethane	0.539	0.513	4.8	83	0.00
19	P	n-Nitroso-di-n-propylamine	1.069	0.943	11.8	79	0.00
20		3+4-Methylphenols	1.533	1.366	10.9	83	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	89	0.00
22		Acetophenone	0.519	0.468	9.8	80	0.00
23	S	Nitrobenzene-d5	0.361	0.358	0.8	82	0.00
24		Nitrobenzene	0.408	0.398	2.5	82	0.00
25		Isophorone	0.787	0.696	11.6	77	0.00
26	C	2-Nitrophenol	0.150	0.161	-7.3	82	0.00
27		2,4-Dimethylphenol	0.303	0.283	6.6	80	0.00
28		bis(2-Chloroethoxy)methane	0.421	0.380	9.7	80	0.00
29	C	2,4-Dichlorophenol	0.281	0.262	6.8	79	0.00
30		1,2,4-Trichlorobenzene	0.297	0.275	7.4	81	0.00
31		Naphthalene	1.004	0.919	8.5	81	0.00
32		Benzoic acid	0.202	0.202	0.0	86	-0.02
33		4-Chloroaniline	0.407	0.368	9.6	78	0.00
34	C	Hexachlorobutadiene	0.169	0.160	5.3	83	0.00
35		Caprolactam	0.091	0.078	14.3	70	-0.01
36	C	4-Chloro-3-methylphenol	0.328	0.290	11.6	75	0.00
37		2-Methylnaphthalene	0.657	0.596	9.3	79	0.00
38		1-Methylnaphthalene	0.622	0.560	10.0	78	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	82	0.00
40		1,2,4,5-Tetrachlorobenzene	0.564	0.539	4.4	79	0.00
41	P	Hexachlorocyclopentadiene	0.280	0.288	-2.9	80	0.00
42	S	2,4,6-Tribromophenol	0.187	0.181	3.2	73	0.00
43	C	2,4,6-Trichlorophenol	0.406	0.398	2.0	77	0.00
44		2,4,5-Trichlorophenol	0.432	0.413	4.4	75	0.00
45	S	2-Fluorobiphenyl	1.501	1.275	15.1	80	0.00
46		1,1'-Biphenyl	1.676	1.572	6.2	77	0.00
47		2-Chloronaphthalene	1.282	1.209	5.7	78	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081721\
Data File : BF125073.D
Acq On : 17 Aug 2021 17:30
Operator : JU/CG
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF081721

Quant Time: Aug 17 18:18:39 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 17 16:45:36 2021
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	0.413	0.435	-5.3	75	0.00
49		Acenaphthylene	2.074	1.947	6.1	77	0.00
50		Dimethylphthalate	1.377	1.259	8.6	73	0.00
51		2,6-Dinitrotoluene	0.273	0.273	0.0	72	0.00
52	C	Acenaphthene	1.260	1.183	6.1	77	0.00
53		3-Nitroaniline	0.340	0.339	0.3	72	0.00
54	P	2,4-Dinitrophenol	0.110	0.107	2.7	73	0.00
55		Dibenzofuran	1.845	1.722	6.7	76	0.00
56	P	4-Nitrophenol	0.275	0.272	1.1	70	0.00
57		2,4-Dinitrotoluene	0.325	0.342	-5.2	72	0.00
58		Fluorene	1.342	1.213	9.6	75	0.00
59		2,3,4,6-Tetrachlorophenol	0.313	0.305	2.6	75	0.00
60		Diethylphthalate	1.469	1.346	8.4	73	0.00
61		4-Chlorophenyl-phenylether	0.608	0.563	7.4	76	0.00
62		4-Nitroaniline	0.308	0.289	6.2	66	-0.01
63		Azobenzene	1.668	1.545	7.4	74	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	78	0.00
65		4,6-Dinitro-2-methylphenol	0.087	0.091	-4.6	71	0.00
66	c	n-Nitrosodiphenylamine	0.713	0.674	5.5	73	0.00
67		4-Bromophenyl-phenylether	0.211	0.207	1.9	73	0.00
68		Hexachlorobenzene	0.232	0.225	3.0	73	0.00
69		Atrazine	0.188	0.174	7.4	70	0.00
70	C	Pentachlorophenol	0.135	0.134	0.7	68	0.00
71		Phenanthrene	1.072	1.008	6.0	72	0.00
72		Anthracene	1.068	0.999	6.5	71	0.00
73		Carbazole	1.105	1.010	8.6	69	0.00
74		Di-n-butylphthalate	1.243	1.182	4.9	70	0.00
75	C	Fluoranthene	1.067	1.002	6.1	69	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	75	0.00
77		Benzidine	0.860	0.730	15.1	63	0.00
78		Pyrene	2.003	1.767	11.8	68	0.00
79	S	Terphenyl-d14	1.389	1.123	19.2	68	0.00
80		Butylbenzylphthalate	0.806	0.735	8.8	65	0.00
81		Benzo(a)anthracene	1.271	1.219	4.1	68	0.00
82		3,3'-Dichlorobenzidine	0.427	0.396	7.3	68	0.00
83		Chrysene	1.290	1.225	5.0	69	0.00
84		Bis(2-ethylhexyl)phthalate	0.957	0.923	3.6	68	0.00
85	c	Di-n-octyl phthalate	1.386	1.506	-8.7	69	0.00
86	I	Perylene-d12	1.000	1.000	0.0	93	0.00
87		Indeno(1,2,3-cd)pyrene	0.693	0.842	-21.5	109	0.00
88		Benzo(b)fluoranthene	1.357	1.175	13.4	73	0.00
89		Benzo(k)fluoranthene	1.285	1.189	7.5	81	0.00
90	C	Benzo(a)pyrene	1.137	1.036	8.9	78	0.00
91		Dibenzo(a,h)anthracene	0.658	0.797	-21.1	108	0.00
92		Benzo(g,h,i)perylene	0.630	0.680	-7.9	110	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF081721\
Data File : BF125073.D
Acq On : 17 Aug 2021 17:30
Operator : JU/CG
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
ICVBF081721

Quant Time: Aug 17 18:18:39 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 17 16:45:36 2021
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