Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP022122\
Data File : BP009229.D
Acq On : 21 Feb 2022 11:54
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Feb 22 01:49:13 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP020722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Feb 19 00:46:08 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	86	0.00
2		1,4-Dioxane	0.369	0.411	-11.4	98	0.00
3		Pyridine	1.079	1.148	-6.4	91	0.00
4		n-Nitrosodimethylamine	0.413	0.397	3.9	79	0.00
5	S	2-Fluorophenol	1.121	1.092	2.6	85	0.00
6		Aniline	1.677	1.650	1.6	87	0.00
7	S	Phenol-d6	1.477	1.447	2.0	86	0.00
8		2-Chlorophenol	1.265	1.224	3.2	86	0.00
9		Benzaldehyde	0.748	0.680	9.1	83	0.00
10	C	Phenol	1.485	1.430	3.7	85	0.00
11		bis(2-Chloroethyl)ether	1.153	1.146	0.6	89	0.00
12		1,3-Dichlorobenzene	1.463	1.418	3.1	87	0.00
13	C	1,4-Dichlorobenzene	1.494	1.448	3.1	86	0.00
14		1,2-Dichlorobenzene	1.448	1.398	3.5	86	0.00
15		Benzyl Alcohol	0.940	1.097	-16.7	102	0.00
16		2,2'-oxybis(1-Chloropropane	1.300	1.251	3.8	88	0.00
17		2-Methylphenol	0.996	0.991	0.5	89	0.00
18		Hexachloroethane	0.495	0.481	2.8	87	0.00
19	P	n-Nitroso-di-n-propylamine	0.846	0.915	-8.2	97	0.00
20		3+4-Methylphenols	1.363	1.386	-1.7	90	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	93	0.00
22		Acetophenone	0.468	0.465	0.6	92	0.00
23	S	Nitrobenzene-d5	0.355	0.345	2.8	90	0.00
24		Nitrobenzene	0.335	0.320	4.5	89	0.00
25		Isophorone	0.589	0.602	-2.2	99	0.00
26	C	2-Nitrophenol	0.173	0.178	-2.9	94	0.00
27		2,4-Dimethylphenol	0.259	0.256	1.2	93	0.00
28		bis(2-Chloroethoxy)methane	0.400	0.394	1.5	95	0.00
29	C	2,4-Dichlorophenol	0.324	0.321	0.9	92	0.00
30		1,2,4-Trichlorobenzene	0.381	0.379	0.5	94	0.00
31		Naphthalene	1.020	0.984	3.5	92	0.00
32		Benzoic acid	0.202	0.132	34.7#	65	-0.01
33		4-Chloroaniline	0.412	0.418	-1.5	97	0.00
34	C	Hexachlorobutadiene	0.234	0.238	-1.7	96	0.00
35		Caprolactam	0.095	0.109	-14.7	117	0.00
36	C	4-Chloro-3-methylphenol	0.309	0.319	-3.2	102	0.00
37		2-Methylnaphthalene	0.722	0.721	0.1	98	0.00
38		1-Methylnaphthalene	0.705	0.708	-0.4	98	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	112	0.00
40		1,2,4,5-Tetrachlorobenzene	0.637	0.592	7.1	102	0.00
41	P	Hexachlorocyclopentadiene	0.186	0.112	39.8#	66	0.00
42	S	2,4,6-Tribromophenol	0.267	0.282	-5.6	125	0.00
43	C	2,4,6-Trichlorophenol	0.422	0.392	7.1	101	0.00
44		2,4,5-Trichlorophenol	0.453	0.414	8.6	101	0.00
45	S	2-Fluorobiphenyl	1.429	1.336	6.5	103	0.00
46		1,1'-Biphenyl	1.469	1.349	8.2	102	0.00
47		2-Chloronaphthalene	1.168	1.082	7.4	103	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.271	0.273	-0.7	113	0.00
49		Acenaphthylene	1.760	1.664	5.5	106	0.00
50		Dimethylphthalate	1.426	1.410	1.1	117	0.00
51		2,6-Dinitrotoluene	0.297	0.308	-3.7	120	0.00
52	C	Acenaphthene	1.071	1.008	5.9	107	0.00
53		3-Nitroaniline	0.304	0.313	-3.0	117	0.00
54	P	2,4-Dinitrophenol	0.157	0.157	0.0	117	0.01
55		Dibenzofuran	1.805	1.747	3.2	112	0.00
56	P	4-Nitrophenol	0.218	0.340	-56.0#	184#	0.01
57		2,4-Dinitrotoluene	0.388	0.433	-11.6	129	0.00
58		Fluorene	1.386	1.386	0.0	116	0.00
59		2,3,4,6-Tetrachlorophenol	0.396	0.393	0.8	115	0.00
60		Diethylphthalate	1.348	1.377	-2.2	124	0.00
61		4-Chlorophenyl-phenylether	0.753	0.759	-0.8	118	0.00
62		4-Nitroaniline	0.306	0.309	-1.0	115	0.00
63		Azobenzene	1.166	1.140	2.2	114	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	134	0.00
65		4,6-Dinitro-2-methylphenol	0.123	0.116	5.7	129	0.00
66	c	n-Nitrosodiphenylamine	0.607	0.549	9.6	120	0.00
67		4-Bromophenyl-phenylether	0.232	0.217	6.5	126	0.00
68		Hexachlorobenzene	0.260	0.247	5.0	129	0.00
69		Atrazine	0.201	0.202	-0.5	137	0.00
70	C	Pentachlorophenol	0.137	0.119	13.1	116	0.01
71		Phenanthrene	1.083	1.011	6.6	126	0.00
72		Anthracene	1.060	1.007	5.0	127	0.00
73		Carbazole	1.016	0.984	3.1	130	0.00
74		Di-n-butylphthalate	1.029	1.062	-3.2	143	0.00
75	C	Fluoranthene	1.214	1.267	-4.4	140	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	143	0.00
77		Benzidine	0.418	0.363	13.2	104	0.00
78		Pyrene	1.423	1.212	14.8	136	0.00
79	S	Terphenyl-d14	1.112	0.976	12.2	140	0.00
80		Butylbenzylphthalate	0.489	0.479	2.0	154#	0.00
81		Benzo(a)anthracene	1.289	1.262	2.1	141	0.00
82		3,3'-Dichlorobenzidine	0.447	0.436	2.5	131	0.00
83		Chrysene	1.242	1.176	5.3	138	0.00
84		Bis(2-ethylhexyl)phthalate	0.647	0.682	-5.4	162#	0.00
85	c	Di-n-octyl phthalate	1.035	1.140	-10.1	152#	0.00
86	I	Perylene-d12	1.000	1.000	0.0	128	0.00
87		Indeno(1,2,3-cd)pyrene	1.486	1.266	14.8	100	0.00
88		Benzo(b)fluoranthene	1.231	1.213	1.5	129	0.00
89		Benzo(k)fluoranthene	1.228	1.182	3.7	126	0.00
90	C	Benzo(a)pyrene	1.027	0.998	2.8	123	0.00
91		Dibenzo(a,h)anthracene	1.221	1.052	13.8	101	-0.01
92		Benzo(g,h,i)perylene	1.216	0.964	20.7	93	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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