Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP070721\
Data File : BP006134.D
Acq On : 07 Jul 2021 16:00
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : ICVBP070721

Quant Time: Jul 07 16:33:02 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP070721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 07 16:08:13 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	85	0.00
2		1,4-Dioxane	0.456	0.486	-6.6	93	0.00
3		Pyridine	1.159	1.304	-12.5	94	0.00
4		n-Nitrosodimethylamine	0.456	0.505	-10.7	97	0.00
5	S	2-Fluorophenol	1.162	1.234	-6.2	90	-0.01
6		Aniline	1.622	1.695	-4.5	89	0.00
7	S	Phenol-d6	1.537	1.604	-4.4	88	0.00
8		2-Chlorophenol	1.361	1.399	-2.8	87	0.00
9		Benzaldehyde	0.839	0.879	-4.8	89	0.00
10	C	Phenol	1.477	1.555	-5.3	90	-0.01
11		bis(2-Chloroethyl)ether	1.116	1.167	-4.6	88	0.00
12		1,3-Dichlorobenzene	1.459	1.472	-0.9	86	0.00
13	C	1,4-Dichlorobenzene	1.490	1.478	0.8	84	0.00
14		1,2-Dichlorobenzene	1.418	1.406	0.8	84	0.00
15		Benzyl Alcohol	0.930	1.004	-8.0	90	0.00
16		2,2'-oxybis(1-Chloropropane	1.325	1.350	-1.9	89	0.00
17		2-Methylphenol	1.005	1.026	-2.1	86	0.00
18		Hexachloroethane	0.496	0.482	2.8	83	0.00
19	P	n-Nitroso-di-n-propylamine	0.838	0.852	-1.7	87	0.00
20		3+4-Methylphenols	1.357	1.363	-0.4	84	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	82	0.00
22		Acetophenone	0.455	0.475	-4.4	85	0.00
23	S	Nitrobenzene-d5	0.317	0.334	-5.4	86	0.00
24		Nitrobenzene	0.320	0.339	-5.9	87	0.00
25		Isophorone	0.601	0.631	-5.0	85	0.00
26	C	2-Nitrophenol	0.178	0.185	-3.9	80	0.00
27		2,4-Dimethylphenol	0.263	0.273	-3.8	82	0.00
28		bis(2-Chloroethoxy)methane	0.333	0.343	-3.0	83	0.00
29	C	2,4-Dichlorophenol	0.289	0.298	-3.1	81	0.00
30		1,2,4-Trichlorobenzene	0.307	0.306	0.3	79	0.00
31		Naphthalene	1.043	1.053	-1.0	83	0.00
32		Benzoic acid	0.177	0.198	-11.9	87	0.00
33		4-Chloroaniline	0.402	0.419	-4.2	82	0.00
34	C	Hexachlorobutadiene	0.169	0.167	1.2	77	0.00
35		Caprolactam	0.091	0.098	-7.7	85	0.00
36	C	4-Chloro-3-methylphenol	0.312	0.330	-5.8	85	0.00
37		2-Methylnaphthalene	0.723	0.735	-1.7	82	0.00
38		1-Methylnaphthalene	0.685	0.698	-1.9	82	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	82	0.00
40		1,2,4,5-Tetrachlorobenzene	0.496	0.492	0.8	78	0.00
41	P	Hexachlorocyclopentadiene	0.076	0.080	-5.3	84	0.00
42	S	2,4,6-Tribromophenol	0.238	0.228	4.2	72	0.00
43	C	2,4,6-Trichlorophenol	0.359	0.372	-3.6	82	0.00
44		2,4,5-Trichlorophenol	0.385	0.402	-4.4	82	0.00
45	S	2-Fluorobiphenyl	1.353	1.366	-1.0	82	0.00
46		1,1'-Biphenyl	1.502	1.533	-2.1	84	0.00
47		2-Chloronaphthalene	1.155	1.185	-2.6	83	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.287	0.316	-10.1	89	0.00
49		Acenaphthylene	1.860	1.909	-2.6	83	0.00
50		Dimethylphthalate	1.438	1.470	-2.2	83	0.00
51		2,6-Dinitrotoluene	0.324	0.341	-5.2	84	0.00
52	C	Acenaphthene	1.131	1.139	-0.7	83	0.00
53		3-Nitroaniline	0.328	0.355	-8.2	85	0.00
54	P	2,4-Dinitrophenol	0.132	0.150	-13.6	87	0.00
55		Dibenzofuran	1.761	1.784	-1.3	83	0.00
56	P	4-Nitrophenol	0.209	0.234	-12.0	89	0.00
57		2,4-Dinitrotoluene	0.437	0.463	-5.9	83	0.00
58		Fluorene	1.415	1.414	0.1	82	0.00
59		2,3,4,6-Tetrachlorophenol	0.301	0.308	-2.3	79	0.00
60		Diethylphthalate	1.475	1.488	-0.9	83	0.00
61		4-Chlorophenyl-phenylether	0.620	0.610	1.6	79	0.00
62		4-Nitroaniline	0.302	0.331	-9.6	84	0.00
63		Azobenzene	1.231	1.283	-4.2	87	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	79	0.00
65		4,6-Dinitro-2-methylphenol	0.121	0.131	-8.3	81	0.00
66	c	n-Nitrosodiphenylamine	0.651	0.670	-2.9	80	0.00
67		4-Bromophenyl-phenylether	0.197	0.193	2.0	74	0.00
68		Hexachlorobenzene	0.234	0.226	3.4	72	0.00
69		Atrazine	0.204	0.211	-3.4	80	0.00
70	C	Pentachlorophenol	0.108	0.110	-1.9	76	0.00
71		Phenanthrene	1.130	1.142	-1.1	80	0.00
72		Anthracene	1.107	1.123	-1.4	79	0.00
73		Carbazole	1.097	1.133	-3.3	81	0.00
74		Di-n-butylphthalate	1.320	1.384	-4.8	82	0.00
75	C	Fluoranthene	1.238	1.237	0.1	78	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	77	0.00
77		Benzidine	0.593	0.675	-13.8	81	0.00
78		Pyrene	1.352	1.367	-1.1	77	0.00
79	S	Terphenyl-d14	1.014	1.026	-1.2	75	0.01
80		Butylbenzylphthalate	0.671	0.710	-5.8	82	0.01
81		Benzo(a)anthracene	1.265	1.289	-1.9	78	0.01
82		3,3'-Dichlorobenzidine	0.371	0.376	-1.3	75	0.00
83		Chrysene	1.250	1.269	-1.5	78	0.01
84		Bis(2-ethylhexyl)phthalate	0.990	1.059	-7.0	85	0.00
85	c	Di-n-octyl phthalate	1.743	1.894	-8.7	86	0.01
86	I	Perylene-d12	1.000	1.000	0.0	83	0.02
87		Indeno(1,2,3-cd)pyrene	1.475	1.502	-1.8	82	0.02
88		Benzo(b)fluoranthene	1.213	1.233	-1.6	80	0.02
89		Benzo(k)fluoranthene	1.186	1.207	-1.8	85	0.01
90	C	Benzo(a)pyrene	1.138	1.153	-1.3	83	0.02
91		Dibenzo(a,h)anthracene	1.270	1.298	-2.2	82	0.03
92		Benzo(g,h,i)perylene	1.250	1.266	-1.3	83	0.03

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Compound	AvgRF	CCRF	%Dev Area%	Dev(min)

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