

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG013024\
 Data File : BG060478.D
 Acq On : 30 Jan 2024 11:06
 Operator : MA/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Jan 30 11:46:51 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG010824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 09 04:07:12 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	88	-0.01
2	1,4-Dioxane	0.555	0.588	-5.9	81	0.00
3	Pyridine	1.567	1.510	3.6	71	0.00
4	n-Nitrosodimethylamine	0.490	0.410	16.3	65	0.00
5 S	2-Fluorophenol	1.216	1.164	4.3	90	0.00
6	Aniline	1.800	1.792	0.4	81	0.00
7 S	Phenol-d6	1.801	1.807	-0.3	76	0.00
8	2-Chlorophenol	1.205	1.243	-3.2	88	0.00
9	Benzaldehyde	1.034	0.871	15.8	68	0.00
10 C	Phenol	1.805	1.804	0.1	75	0.00
11	bis(2-Chloroethyl)ether	1.471	1.392	5.4	82	0.00
12	1,3-Dichlorobenzene	1.513	1.518	-0.3	87	0.00
13 C	1,4-Dichlorobenzene	1.535	1.534	0.1	89	0.00
14	1,2-Dichlorobenzene	1.578	1.484	6.0	74	0.00
15	Benzyl Alcohol	1.166	1.237	-6.1	87	0.00
16	2,2'-oxybis(1-Chloropropane	1.785	1.798	-0.7	79	0.00
17	2-Methylphenol	1.240	1.255	-1.2	77	0.00
18	Hexachloroethane	0.538	0.529	1.7	77	0.00
19 P	n-Nitroso-di-n-propylamine	1.246	1.269	-1.8	75	0.00
20	3+4-Methylphenols	1.672	1.789	-7.0	80	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	80	0.00
22	Acetophenone	0.550	0.548	0.4	79	0.00
23 S	Nitrobenzene-d5	0.424	0.439	-3.5	83	0.00
24	Nitrobenzene	0.439	0.422	3.9	76	0.00
25	Isophorone	0.851	0.846	0.6	79	0.00
26 C	2-Nitrophenol	0.161	0.182	-13.0	88	0.00
27	2,4-Dimethylphenol	0.213	0.226	-6.1	83	0.00
28	bis(2-Chloroethoxy)methane	0.520	0.495	4.8	76	0.00
29 C	2,4-Dichlorophenol	0.346	0.357	-3.2	81	0.00
30	1,2,4-Trichlorobenzene	0.430	0.434	-0.9	80	0.00
31	Naphthalene	1.048	1.041	0.7	81	0.00
32	Benzoic acid	0.154	0.185	-20.1	88	0.00
33	4-Chloroaniline	0.404	0.410	-1.5	82	0.00
34 C	Hexachlorobutadiene	0.280	0.292	-4.3	89	0.00
35	Caprolactam	0.111	0.123	-10.8	89	0.00
36 C	4-Chloro-3-methylphenol	0.351	0.381	-8.5	87	0.00
37	2-Methylnaphthalene	0.778	0.789	-1.4	84	0.00
38	1-Methylnaphthalene	0.781	0.785	-0.5	84	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	83	0.00
40	1,2,4,5-Tetrachlorobenzene	0.710	0.709	0.1	87	0.00
41 P	Hexachlorocyclopentadiene	0.236	0.261	-10.6	89	0.00
42 S	2,4,6-Tribromophenol	0.294	0.318	-8.2	90	0.00
43 C	2,4,6-Trichlorophenol	0.452	0.473	-4.6	88	0.00
44	2,4,5-Trichlorophenol	0.499	0.537	-7.6	90	0.00
45 S	2-Fluorobiphenyl	1.439	1.433	0.4	89	0.00
46	1,1'-Biphenyl	1.510	1.467	2.8	84	0.00
47	2-Chloronaphthalene	1.293	1.249	3.4	83	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG013024\
 Data File : BG060478.D
 Acq On : 30 Jan 2024 11:06
 Operator : MA/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Jan 30 11:46:51 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG010824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 09 04:07:12 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	0.322	0.348	-8.1	88	0.00
49	Acenaphthylene	1.781	1.783	-0.1	84	0.00
50	Dimethylphthalate	1.528	1.506	1.4	84	0.00
51	2,6-Dinitrotoluene	0.326	0.358	-9.8	88	0.00
52 C	Acenaphthene	1.109	1.096	1.2	84	0.00
53	3-Nitroaniline	0.297	0.323	-8.8	88	0.00
54 P	2,4-Dinitrophenol	0.155	0.195	-25.8#	98	0.00
55	Dibenzofuran	1.851	1.837	0.8	84	0.00
56 P	4-Nitrophenol	0.220	0.252	-14.5	86	0.00
57	2,4-Dinitrotoluene	0.447	0.509	-13.9	93	0.00
58	Fluorene	1.533	1.521	0.8	85	0.00
59	2,3,4,6-Tetrachlorophenol	0.423	0.473	-11.8	93	0.00
60	Diethylphthalate	1.467	1.494	-1.8	86	0.00
61	4-Chlorophenyl-phenylether	0.845	0.861	-1.9	89	0.00
62	4-Nitroaniline	0.316	0.344	-8.9	89	0.00
63	Azobenzene	1.484	1.416	4.6	81	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	87	0.00
65	4,6-Dinitro-2-methylphenol	0.111	0.127	-14.4	98	0.00
66 c	n-Nitrosodiphenylamine	0.531	0.529	0.4	86	0.00
67	4-Bromophenyl-phenylether	0.220	0.226	-2.7	89	0.00
68	Hexachlorobenzene	0.260	0.266	-2.3	90	0.00
69	Atrazine	0.200	0.186	7.0	82	0.00
70 C	Pentachlorophenol	0.140	0.170	-21.4#	98	0.00
71	Phenanthrene	0.967	0.966	0.1	88	0.00
72	Anthracene	0.965	0.970	-0.5	88	0.00
73	Carbazole	0.910	0.891	2.1	85	0.00
74	Di-n-butylphthalate	0.936	0.952	-1.7	87	0.00
75 C	Fluoranthene	1.161	1.149	1.0	89	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	87	0.00
77	Benzydine	0.435	0.368	15.4	72	0.00
78	Pyrene	1.277	1.240	2.9	89	0.00
79 S	Terphenyl-d14	0.938	0.848	9.6	95	0.00
80	Butylbenzylphthalate	0.439	0.468	-6.6	91	0.00
81	Benzo(a)anthracene	1.238	1.237	0.1	89	0.00
82	3,3'-Dichlorobenzidine	0.474	0.474	0.0	86	0.00
83	Chrysene	1.218	1.199	1.6	89	0.00
84	Bis(2-ethylhexyl)phthalate	0.664	0.678	-2.1	87	0.00
85 c	Di-n-octyl phthalate	1.076	1.123	-4.4	89	0.00
86 I	Perylene-d12	1.000	1.000	0.0	87	0.00
87	Indeno(1,2,3-cd)pyrene	1.391	1.375	1.2	87	0.00
88	Benzo(b)fluoranthene	1.182	1.185	-0.3	86	0.00
89	Benzo(k)fluoranthene	1.166	1.131	3.0	83	0.00
90 C	Benzo(a)pyrene	1.067	1.058	0.8	87	0.00
91	Dibenzo(a,h)anthracene	1.136	1.121	1.3	85	0.00
92	Benzo(g,h,i)perylene	1.164	1.145	1.6	86	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG013024\
Data File : BG060478.D
Acq On : 30 Jan 2024 11:06
Operator : MA/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Jan 30 11:46:51 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG010824.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 09 04:07:12 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 = 1