

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG031924\
 Data File : BG060677.D
 Acq On : 19 Mar 2024 11:01
 Operator : MA/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Mar 19 11:46:34 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 18 14:26:31 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	85	0.00
2	1,4-Dioxane	0.533	0.667	-25.1#	109	0.00
3	Pyridine	1.574	1.924	-22.2	103	0.00
4	n-Nitrosodimethylamine	0.772	1.017	-31.7#	111	0.00
5 S	2-Fluorophenol	1.272	1.468	-15.4	96	0.00
6	Aniline	2.255	2.175	3.5	79	0.00
7 S	Phenol-d6	1.846	1.760	4.7	78	0.00
8	2-Chlorophenol	1.391	1.315	5.5	78	0.00
9	Benzaldehyde	1.033	0.960	7.1	75	0.00
10 C	Phenol	1.852	1.746	5.7	77	0.00
11	bis(2-Chloroethyl)ether	1.476	1.385	6.2	78	0.00
12	1,3-Dichlorobenzene	1.500	1.396	6.9	79	0.00
13 C	1,4-Dichlorobenzene	1.520	1.404	7.6	77	0.00
14	1,2-Dichlorobenzene	1.510	1.394	7.7	78	0.00
15	Benzyl Alcohol	1.459	1.341	8.1	74	0.00
16	2,2'-oxybis(1-Chloropropane	2.866	2.689	6.2	78	0.00
17	2-Methylphenol	1.422	1.368	3.8	78	0.00
18	Hexachloroethane	0.519	0.485	6.6	78	0.00
19 P	n-Nitroso-di-n-propylamine	1.279	1.259	1.6	80	0.00
20	3+4-Methylphenols	1.918	1.854	3.3	78	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	82	0.00
22	Acetophenone	0.536	0.525	2.1	78	0.00
23 S	Nitrobenzene-d5	0.384	0.351	8.6	72	0.00
24	Nitrobenzene	0.384	0.367	4.4	75	0.00
25	Isophorone	0.769	0.767	0.3	80	0.00
26 C	2-Nitrophenol	0.142	0.121	14.8	65	0.00
27	2,4-Dimethylphenol	0.340	0.318	6.5	77	0.00
28	bis(2-Chloroethoxy)methane	0.461	0.449	2.6	79	0.00
29 C	2,4-Dichlorophenol	0.306	0.300	2.0	78	0.00
30	1,2,4-Trichlorobenzene	0.324	0.304	6.2	77	0.00
31	Naphthalene	1.117	1.067	4.5	78	0.00
32	Benzoic acid	0.199	0.189	5.0	71	0.00
33	4-Chloroaniline	0.471	0.467	0.8	79	0.00
34 C	Hexachlorobutadiene	0.221	0.201	9.0	74	0.00
35	Caprolactam	0.103	0.110	-6.8	85	0.00
36 C	4-Chloro-3-methylphenol	0.374	0.381	-1.9	81	0.00
37	2-Methylnaphthalene	0.772	0.740	4.1	78	0.00
38	1-Methylnaphthalene	0.724	0.705	2.6	79	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	86	0.00
40	1,2,4,5-Tetrachlorobenzene	0.614	0.547	10.9	76	0.00
41 P	Hexachlorocyclopentadiene	0.302	0.250	17.2	68	0.00
42 S	2,4,6-Tribromophenol	0.232	0.245	-5.6	85	0.00
43 C	2,4,6-Trichlorophenol	0.394	0.367	6.9	78	0.00
44	2,4,5-Trichlorophenol	0.467	0.443	5.1	78	0.00
45 S	2-Fluorobiphenyl	1.479	1.359	8.1	79	0.00
46	1,1'-Biphenyl	1.520	1.403	7.7	79	0.00
47	2-Chloronaphthalene	1.147	1.062	7.4	78	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG031924\
 Data File : BG060677.D
 Acq On : 19 Mar 2024 11:01
 Operator : MA/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Mar 19 11:46:34 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 18 14:26:31 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	0.379	0.339	10.6	71	0.00
49	Acenaphthylene	1.869	1.783	4.6	81	0.00
50	Dimethylphthalate	1.500	1.454	3.1	82	0.00
51	2,6-Dinitrotoluene	0.276	0.248	10.1	72	0.00
52 C	Acenaphthene	1.117	1.067	4.5	82	0.00
53	3-Nitroaniline	0.318	0.306	3.8	79	0.00
54 P	2,4-Dinitrophenol	0.118	0.103	12.7	72	0.00
55	Dibenzofuran	1.808	1.726	4.5	82	0.00
56 P	4-Nitrophenol	0.237	0.238	-0.4	83	0.00
57	2,4-Dinitrotoluene	0.348	0.333	4.3	76	0.00
58	Fluorene	1.447	1.415	2.2	83	0.00
59	2,3,4,6-Tetrachlorophenol	0.389	0.383	1.5	81	0.00
60	Diethylphthalate	1.543	1.524	1.2	84	0.00
61	4-Chlorophenyl-phenylether	0.733	0.693	5.5	81	0.00
62	4-Nitroaniline	0.326	0.340	-4.3	84	0.00
63	Azobenzene	1.597	1.573	1.5	83	0.00
64 I	Phenanthrene-d10	1.000	1.000	0.0	88	0.00
65	4,6-Dinitro-2-methylphenol	0.083	0.074	10.8	75	0.00
66 c	n-Nitrosodiphenylamine	0.626	0.591	5.6	83	0.00
67	4-Bromophenyl-phenylether	0.216	0.213	1.4	89	0.00
68	Hexachlorobenzene	0.248	0.231	6.9	83	0.00
69	Atrazine	0.209	0.205	1.9	85	0.00
70 C	Pentachlorophenol	0.161	0.160	0.6	83	0.00
71	Phenanthrene	1.077	1.028	4.5	85	0.00
72	Anthracene	1.089	1.063	2.4	86	0.00
73	Carbazole	0.947	0.942	0.5	87	0.00
74	Di-n-butylphthalate	1.154	1.162	-0.7	86	0.00
75 C	Fluoranthene	1.192	1.190	0.2	87	0.00
76 I	Chrysene-d12	1.000	1.000	0.0	92	0.00
77	Benzidine	0.489	0.602	-23.1	101	0.00
78	Pyrene	1.396	1.337	4.2	87	0.00
79 S	Terphenyl-d14	1.106	1.041	5.9	87	0.00
80	Butylbenzylphthalate	0.528	0.539	-2.1	89	0.00
81	Benzo(a)anthracene	1.377	1.322	4.0	86	0.00
82	3,3'-Dichlorobenzidine	0.464	0.474	-2.2	88	0.00
83	Chrysene	1.337	1.268	5.2	86	0.00
84	Bis(2-ethylhexyl)phthalate	0.779	0.793	-1.8	88	0.00
85 c	Di-n-octyl phthalate	1.283	1.358	-5.8	91	0.00
86 I	Perylene-d12	1.000	1.000	0.0	94	-0.01
87	Indeno(1,2,3-cd)pyrene	1.364	1.338	1.9	90	-0.02
88	Benzo(b)fluoranthene	1.100	1.068	2.9	88	-0.01
89	Benzo(k)fluoranthene	1.156	1.116	3.5	90	0.00
90 C	Benzo(a)pyrene	1.079	1.058	1.9	90	0.00
91	Dibenzo(a,h)anthracene	1.158	1.106	4.5	88	-0.01
92	Benzo(g,h,i)perylene	1.121	1.097	2.1	89	-0.02

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG031924\
Data File : BG060677.D
Acq On : 19 Mar 2024 11:01
Operator : MA/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Mar 19 11:46:34 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG031324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 18 14:26:31 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 = 0