

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090424\
 Data File : BG062656.D
 Acq On : 5 Sep 2024 00:25
 Operator : MA/JU
 Sample : SSTDCCC040
 Misc : LOD-MDL-WATER-8 ng
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Sep 05 04:55:29 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG083024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 31 02:36:01 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	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	111	0.00
2	1,4-Dioxane	40.000	39.293	1.8	111	0.00
3	Pyridine	40.000	40.892	-2.2	112	0.00
4	n-Nitrosodimethylamine	40.000	42.408	-6.0	114	0.00
5 S	2-Fluorophenol	80.000	81.580	-2.0	110	0.00
6	Aniline	40.000	39.358	1.6	104	0.00
7 S	Phenol-d6	80.000	82.910	-3.6	110	0.00
8	2-Chlorophenol	40.000	40.430	-1.1	110	0.00
9	Benzaldehyde	40.000	45.589	-14.0	128	0.00
10 C	Phenol	40.000	41.394	-3.5	110	0.00
11	bis(2-Chloroethyl)ether	40.000	39.851	0.4	108	0.00
12	1,3-Dichlorobenzene	40.000	40.662	-1.7	112	0.00
13 C	1,4-Dichlorobenzene	40.000	40.131	-0.3	110	0.00
14	1,2-Dichlorobenzene	40.000	39.857	0.4	108	0.00
15	Benzyl Alcohol	40.000	40.313	-0.8	105	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	36.434	8.9	104	0.00
17	2-Methylphenol	40.000	39.537	1.2	106	0.00
18	Hexachloroethane	40.000	39.451	1.4	108	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	38.092	4.8	100	0.00
20	3+4-Methylphenols	40.000	39.318	1.7	105	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	104	0.00
22	Acetophenone	40.000	40.357	-0.9	103	0.00
23 S	Nitrobenzene-d5	80.000	86.117	-7.6	109	0.00
24	Nitrobenzene	40.000	41.956	-4.9	104	0.00
25	Isophorone	40.000	38.806	3.0	98	0.00
26 C	2-Nitrophenol	40.000	45.350	-13.4	112	0.00
27	2,4-Dimethylphenol	40.000	41.027	-2.6	102	0.00
28	bis(2-Chloroethoxy)methane	40.000	39.692	0.8	98	0.00
29 C	2,4-Dichlorophenol	40.000	42.312	-5.8	106	0.00
30	1,2,4-Trichlorobenzene	40.000	41.822	-4.6	108	0.00
31	Naphthalene	40.000	40.900	-2.2	104	0.00
32	Benzoic acid	40.000	39.225	1.9	97	0.00
33	4-Chloroaniline	40.000	41.480	-3.7	105	0.00
34 C	Hexachlorobutadiene	40.000	42.548	-6.4	112	0.00
35	Caprolactam	40.000	41.914	-4.8	104	0.00
36 C	4-Chloro-3-methylphenol	40.000	40.753	-1.9	101	0.00
37	2-Methylnaphthalene	40.000	40.759	-1.9	103	0.00
38	1-Methylnaphthalene	40.000	40.162	-0.4	101	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	103	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	42.370	-5.9	108	0.00
41 P	Hexachlorocyclopentadiene	40.000	45.432	-13.6	112	0.00
42 S	2,4,6-Tribromophenol	80.000	90.109	-12.6	112	0.00
43 C	2,4,6-Trichlorophenol	40.000	42.170	-5.4	105	0.00
44	2,4,5-Trichlorophenol	40.000	42.981	-7.5	106	0.00
45 S	2-Fluorobiphenyl	80.000	80.573	-0.7	102	0.00
46	1,1'-Biphenyl	40.000	41.040	-2.6	104	0.00
47	2-Chloronaphthalene	40.000	41.397	-3.5	105	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090424\
 Data File : BG062656.D
 Acq On : 5 Sep 2024 00:25
 Operator : MA/JU
 Sample : SSTDCCC040
 Misc : LOD-MDL-WATER-8 ng
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 LabSampleId :
 SSTDCCC040

Quant Time: Sep 05 04:55:29 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG083024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 31 02:36:01 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	Amount	Calc.	%Dev	Area%	Dev(min)
48	2-Nitroaniline	40.000	45.741	-14.4	113	0.00
49	Acenaphthylene	40.000	40.533	-1.3	101	0.00
50	Dimethylphthalate	40.000	39.634	0.9	100	0.00
51	2,6-Dinitrotoluene	40.000	44.799	-12.0	107	0.00
52 C	Acenaphthene	40.000	40.919	-2.3	104	0.00
53	3-Nitroaniline	40.000	45.712	-14.3	109	0.00
54 P	2,4-Dinitrophenol	40.000	47.177	-17.9	117	0.00
55	Dibenzofuran	40.000	41.215	-3.0	105	0.00
56 P	4-Nitrophenol	40.000	45.903	-14.8	110	0.00
57	2,4-Dinitrotoluene	40.000	48.313	-20.8	118	0.00
58	Fluorene	40.000	41.405	-3.5	106	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	43.141	-7.9	111	0.00
60	Diethylphthalate	40.000	40.530	-1.3	104	0.00
61	4-Chlorophenyl-phenylether	40.000	42.241	-5.6	109	0.00
62	4-Nitroaniline	40.000	45.867	-14.7	113	0.00
63	Azobenzene	40.000	40.388	-1.0	102	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	107	0.00
65	4,6-Dinitro-2-methylphenol	40.000	46.464	-16.2	122	0.00
66 c	n-Nitrosodiphenylamine	40.000	40.375	-0.9	106	0.00
67	4-Bromophenyl-phenylether	40.000	42.515	-6.3	112	0.00
68	Hexachlorobenzene	40.000	42.328	-5.8	112	0.00
69	Atrazine	40.000	32.624	18.4	100	0.00
70 C	Pentachlorophenol	40.000	42.345	-5.9	108	0.00
71	Phenanthrene	40.000	40.053	-0.1	105	0.00
72	Anthracene	40.000	40.485	-1.2	105	0.00
73	Carbazole	40.000	40.211	-0.5	107	0.00
74	Di-n-butylphthalate	40.000	40.704	-1.8	106	0.00
75 C	Fluoranthene	40.000	41.398	-3.5	111	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	117	0.00
77	Benzydine	40.000	42.886	-7.2	103	0.00
78	Pyrene	40.000	38.660	3.4	111	0.00
79 S	Terphenyl-d14	80.000	75.575	5.5	112	0.00
80	Butylbenzylphthalate	40.000	40.383	-1.0	114	0.00
81	Benzo(a)anthracene	40.000	39.689	0.8	115	0.00
82	3,3'-Dichlorobenzidine	40.000	40.049	-0.1	115	0.00
83	Chrysene	40.000	40.407	-1.0	117	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	39.650	0.9	112	0.00
85 c	Di-n-octyl phthalate	40.000	38.881	2.8	110	0.00
86 I	Perylene-d12	20.000	20.000	0.0	119	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	40.915	-2.3	120	0.00
88	Benzo(b)fluoranthene	40.000	39.945	0.1	117	0.00
89	Benzo(k)fluoranthene	40.000	40.228	-0.6	117	0.00
90 C	Benzo(a)pyrene	40.000	40.087	-0.2	117	0.00
91	Dibenzo(a,h)anthracene	40.000	40.979	-2.4	121	0.00
92	Benzo(g,h,i)perylene	40.000	41.287	-3.2	120	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG090424\
Data File : BG062656.D
Acq On : 5 Sep 2024 00:25
Operator : MA/JU
Sample : SSTDCCC040
Misc : LOD-MDL-WATER-8 ng
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Sep 05 04:55:29 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG083024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Aug 31 02:36:01 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	Amount	Calc.	%Dev	Area%	Dev(min)
----------	--------	-------	------	-------	----------

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