

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081424\
 Data File : BP021492.D
 Acq On : 14 Aug 2024 09:11
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 LabSampleId :
 SSTDCCC040

Quant Time: Aug 14 10:54:41 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 14 00:00:30 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	99	0.00
2	1,4-Dioxane	40.000	40.008	-0.0	99	0.00
3	Pyridine	40.000	41.223	-3.1	101	0.00
4	n-Nitrosodimethylamine	40.000	41.561	-3.9	104	0.00
5 S	2-Fluorophenol	80.000	80.944	-1.2	101	0.00
6	Aniline	40.000	40.281	-0.7	99	0.00
7 S	Phenol-d6	80.000	81.655	-2.1	101	0.00
8	2-Chlorophenol	40.000	40.969	-2.4	101	0.00
9	Benzaldehyde	40.000	35.474	11.3	93	0.00
10 C	Phenol	40.000	40.468	-1.2	100	0.00
11	bis(2-Chloroethyl)ether	40.000	39.697	0.8	100	0.00
12	1,3-Dichlorobenzene	40.000	39.573	1.1	100	0.00
13 C	1,4-Dichlorobenzene	40.000	39.681	0.8	100	0.00
14	1,2-Dichlorobenzene	40.000	39.347	1.6	100	0.00
15	Benzyl Alcohol	40.000	41.330	-3.3	102	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	39.602	1.0	101	0.00
17	2-Methylphenol	40.000	40.716	-1.8	101	0.00
18	Hexachloroethane	40.000	40.123	-0.3	101	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	40.348	-0.9	102	0.00
20	3+4-Methylphenols	40.000	41.680	-4.2	102	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	102	0.00
22	Acetophenone	40.000	39.808	0.5	101	0.00
23 S	Nitrobenzene-d5	80.000	88.830	-11.0	106	0.00
24	Nitrobenzene	40.000	44.067	-10.2	106	0.00
25	Isophorone	40.000	40.218	-0.5	102	0.00
26 C	2-Nitrophenol	40.000	43.288	-8.2	116	0.00
27	2,4-Dimethylphenol	40.000	40.820	-2.1	104	0.00
28	bis(2-Chloroethoxy)methane	40.000	40.115	-0.3	104	0.00
29 C	2,4-Dichlorophenol	40.000	42.786	-7.0	106	0.00
30	1,2,4-Trichlorobenzene	40.000	39.095	2.3	100	0.00
31	Naphthalene	40.000	39.833	0.4	101	0.00
32	Benzoic acid	40.000	45.836	-14.6	130	0.00
33	4-Chloroaniline	40.000	41.625	-4.1	105	0.00
34 C	Hexachlorobutadiene	40.000	38.887	2.8	99	0.00
35	Caprolactam	40.000	46.333	-15.8	115	0.00
36 C	4-Chloro-3-methylphenol	40.000	43.245	-8.1	106	0.00
37	2-Methylnaphthalene	40.000	40.267	-0.7	103	0.00
38	1-Methylnaphthalene	40.000	40.232	-0.6	103	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	107	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	39.314	1.7	104	0.00
41 P	Hexachlorocyclopentadiene	40.000	39.249	1.9	102	0.00
42 S	2,4,6-Tribromophenol	80.000	85.473	-6.8	108	0.00
43 C	2,4,6-Trichlorophenol	40.000	42.503	-6.3	108	-0.01
44	2,4,5-Trichlorophenol	40.000	43.345	-8.4	110	0.00
45 S	2-Fluorobiphenyl	80.000	78.907	1.4	105	0.00
46	1,1'-Biphenyl	40.000	39.529	1.2	105	0.00
47	2-Chloronaphthalene	40.000	39.429	1.4	105	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081424\
 Data File : BP021492.D
 Acq On : 14 Aug 2024 09:11
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 LabSampleId :
 SSTDCCC040

Quant Time: Aug 14 10:54:41 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081324.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 14 00:00:30 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	43.529	-8.8	125	0.00
49	Acenaphthylene	40.000	40.389	-1.0	107	0.00
50	Dimethylphthalate	40.000	41.603	-4.0	111	-0.01
51	2,6-Dinitrotoluene	40.000	42.366	-5.9	119	0.00
52 C	Acenaphthene	40.000	40.129	-0.3	107	0.00
53	3-Nitroaniline	40.000	44.336	-10.8	123	0.00
54 P	2,4-Dinitrophenol	40.000	45.517	-13.8	134	0.00
55	Dibenzofuran	40.000	40.444	-1.1	108	0.00
56 P	4-Nitrophenol	40.000	45.280	-13.2	127	-0.01
57	2,4-Dinitrotoluene	40.000	44.201	-10.5	126	0.00
58	Fluorene	40.000	40.869	-2.2	109	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	44.934	-12.3	113	0.00
60	Diethylphthalate	40.000	42.041	-5.1	111	-0.01
61	4-Chlorophenyl-phenylether	40.000	40.633	-1.6	109	0.00
62	4-Nitroaniline	40.000	45.820	-14.6	123	0.00
63	Azobenzene	40.000	41.567	-3.9	110	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	113	0.00
65	4,6-Dinitro-2-methylphenol	40.000	43.370	-8.4	134	0.00
66 c	n-Nitrosodiphenylamine	40.000	38.811	3.0	109	-0.01
67	4-Bromophenyl-phenylether	40.000	36.645	8.4	106	0.00
68	Hexachlorobenzene	40.000	38.598	3.5	109	0.00
69	Atrazine	40.000	40.371	-0.9	110	0.00
70 C	Pentachlorophenol	40.000	43.212	-8.0	122	0.00
71	Phenanthrene	40.000	38.900	2.8	110	-0.01
72	Anthracene	40.000	39.780	0.5	112	-0.01
73	Carbazole	40.000	40.782	-2.0	113	0.00
74	Di-n-butylphthalate	40.000	41.262	-3.2	114	0.00
75 C	Fluoranthene	40.000	40.792	-2.0	114	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	117	0.00
77	Benzydine	40.000	38.785	3.0	111	0.00
78	Pyrene	40.000	39.136	2.2	116	0.00
79 S	Terphenyl-d14	80.000	77.133	3.6	116	0.00
80	Butylbenzylphthalate	40.000	38.854	2.9	118	0.00
81	Benzo(a)anthracene	40.000	40.049	-0.1	116	0.00
82	3,3'-Dichlorobenzidine	40.000	42.036	-5.1	116	-0.01
83	Chrysene	40.000	39.619	1.0	115	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	42.105	-5.3	117	0.00
85 c	Di-n-octyl phthalate	40.000	39.410	1.5	118	0.00
86 I	Perylene-d12	20.000	20.000	0.0	117	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	40.343	-0.9	116	-0.01
88	Benzo(b)fluoranthene	40.000	40.651	-1.6	119	0.00
89	Benzo(k)fluoranthene	40.000	39.041	2.4	112	0.00
90 C	Benzo(a)pyrene	40.000	40.278	-0.7	116	0.00
91	Dibenzo(a,h)anthracene	40.000	40.593	-1.5	116	-0.02
92	Benzo(g,h,i)perylene	40.000	40.457	-1.1	116	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081424\
Data File : BP021492.D
Acq On : 14 Aug 2024 09:11
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Aug 14 10:54:41 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP081324.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 14 00:00:30 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