

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021825\
 Data File : BP024080.D
 Acq On : 18 Feb 2025 11:08
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 LabSampleId :
 SSTDCCC040

Quant Time: Feb 18 11:33:26 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP021725.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 18 02:25:07 2025
 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	94	0.00
2	1,4-Dioxane	40.000	39.285	1.8	87	0.00
3	Pyridine	40.000	40.439	-1.1	88	0.00
4	n-Nitrosodimethylamine	40.000	40.343	-0.9	89	0.00
5 S	2-Fluorophenol	80.000	80.436	-0.5	87	0.00
6	Aniline	40.000	40.200	-0.5	90	0.00
7 S	Phenol-d6	80.000	80.912	-1.1	89	0.00
8	2-Chlorophenol	40.000	39.957	0.1	88	0.00
9	Benzaldehyde	40.000	37.984	5.0	87	0.00
10 C	Phenol	40.000	39.724	0.7	89	0.00
11	bis(2-Chloroethyl)ether	40.000	39.085	2.3	88	0.00
12	1,3-Dichlorobenzene	40.000	39.007	2.5	87	0.00
13 C	1,4-Dichlorobenzene	40.000	39.526	1.2	88	0.00
14	1,2-Dichlorobenzene	40.000	39.864	0.3	89	0.00
15	Benzyl Alcohol	40.000	41.607	-4.0	92	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	40.260	-0.6	92	0.00
17	2-Methylphenol	40.000	41.604	-4.0	92	0.00
18	Hexachloroethane	40.000	40.234	-0.6	91	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	42.189	-5.5	94	0.00
20	3+4-Methylphenols	40.000	41.693	-4.2	93	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	98	0.00
22	Acetophenone	40.000	40.022	-0.1	94	0.00
23 S	Nitrobenzene-d5	80.000	80.387	-0.5	92	0.00
24	Nitrobenzene	40.000	39.903	0.2	91	0.00
25	Isophorone	40.000	40.371	-0.9	94	0.00
26 C	2-Nitrophenol	40.000	40.688	-1.7	90	0.00
27	2,4-Dimethylphenol	40.000	39.997	0.0	92	0.00
28	bis(2-Chloroethoxy)methane	40.000	38.673	3.3	91	0.00
29 C	2,4-Dichlorophenol	40.000	39.797	0.5	90	0.00
30	1,2,4-Trichlorobenzene	40.000	38.898	2.8	90	0.00
31	Naphthalene	40.000	38.976	2.6	91	0.00
32	Benzoic acid	40.000	36.269	9.3	92	0.00
33	4-Chloroaniline	40.000	39.327	1.7	92	0.00
34 C	Hexachlorobutadiene	40.000	38.612	3.5	88	0.00
35	Caprolactam	40.000	36.622	8.4	94	0.00
36 C	4-Chloro-3-methylphenol	40.000	40.014	-0.0	92	0.00
37	2-Methylnaphthalene	40.000	38.962	2.6	92	0.00
38	1-Methylnaphthalene	40.000	39.268	1.8	93	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	102	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	39.507	1.2	92	0.00
41 P	Hexachlorocyclopentadiene	40.000	40.217	-0.5	90	0.00
42 S	2,4,6-Tribromophenol	80.000	78.591	1.8	96	0.00
43 C	2,4,6-Trichlorophenol	40.000	39.952	0.1	91	0.00
44	2,4,5-Trichlorophenol	40.000	40.097	-0.2	93	0.01
45 S	2-Fluorobiphenyl	80.000	80.553	-0.7	95	0.00
46	1,1'-Biphenyl	40.000	39.284	1.8	96	0.00
47	2-Chloronaphthalene	40.000	39.873	0.3	96	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021825\
 Data File : BP024080.D
 Acq On : 18 Feb 2025 11:08
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 LabSampleId :
 SSTDCCC040

Quant Time: Feb 18 11:33:26 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP021725.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 18 02:25:07 2025
 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	41.957	-4.9	98	0.00
49	Acenaphthylene	40.000	40.126	-0.3	95	0.01
50	Dimethylphthalate	40.000	38.409	4.0	96	0.00
51	2,6-Dinitrotoluene	40.000	40.995	-2.5	99	0.00
52 C	Acenaphthene	40.000	38.884	2.8	95	0.01
53	3-Nitroaniline	40.000	41.139	-2.8	98	0.01
54 P	2,4-Dinitrophenol	40.000	35.271	11.8	95	0.01
55	Dibenzofuran	40.000	39.210	2.0	96	0.01
56 P	4-Nitrophenol	40.000	39.829	0.4	97	0.01
57	2,4-Dinitrotoluene	40.000	40.980	-2.4	97	0.00
58	Fluorene	40.000	39.670	0.8	96	0.01
59	2,3,4,6-Tetrachlorophenol	40.000	39.489	1.3	96	0.00
60	Diethylphthalate	40.000	38.635	3.4	96	0.01
61	4-Chlorophenyl-phenylether	40.000	39.032	2.4	96	0.02
62	4-Nitroaniline	40.000	37.920	5.2	98	0.00
63	Azobenzene	40.000	39.764	0.6	96	0.02
64 I	Phenanthrene-d10	20.000	20.000	0.0	104	0.01
65	4,6-Dinitro-2-methylphenol	40.000	37.048	7.4	97	0.01
66 c	n-Nitrosodiphenylamine	40.000	39.553	1.1	95	0.01
67	4-Bromophenyl-phenylether	40.000	39.175	2.1	96	0.01
68	Hexachlorobenzene	40.000	38.451	3.9	95	0.02
69	Atrazine	40.000	41.132	-2.8	96	0.00
70 C	Pentachlorophenol	40.000	38.359	4.1	96	0.01
71	Phenanthrene	40.000	38.929	2.7	96	0.02
72	Anthracene	40.000	39.677	0.8	95	0.01
73	Carbazole	40.000	39.772	0.6	96	0.01
74	Di-n-butylphthalate	40.000	40.952	-2.4	100	0.00
75 C	Fluoranthene	40.000	38.780	3.0	97	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	106	0.00
77	Benzydine	40.000	58.129	-45.3#	105	0.00
78	Pyrene	40.000	39.287	1.8	96	0.00
79 S	Terphenyl-d14	80.000	79.945	0.1	99	0.00
80	Butylbenzylphthalate	40.000	40.020	-0.1	102	0.00
81	Benzo(a)anthracene	40.000	39.517	1.2	100	0.00
82	3,3'-Dichlorobenzidine	40.000	43.032	-7.6	103	0.00
83	Chrysene	40.000	38.919	2.7	98	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	40.176	-0.4	101	-0.01
85 c	Di-n-octyl phthalate	40.000	38.226	4.4	103	-0.01
86 I	Perylene-d12	20.000	20.000	0.0	108	-0.02
87	Indeno(1,2,3-cd)pyrene	40.000	40.038	-0.1	97	-0.03
88	Benzo(b)fluoranthene	40.000	38.895	2.8	98	-0.02
89	Benzo(k)fluoranthene	40.000	39.997	0.0	100	-0.01
90 C	Benzo(a)pyrene	40.000	39.767	0.6	98	-0.02
91	Dibenzo(a,h)anthracene	40.000	40.030	-0.1	97	-0.04
92	Benzo(g,h,i)perylene	40.000	39.568	1.1	95	-0.04

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP021825\
Data File : BP024080.D
Acq On : 18 Feb 2025 11:08
Operator : RC/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Feb 18 11:33:26 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP021725.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 18 02:25:07 2025
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