

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111021\
 Data File : BP007921.D
 Acq On : 10 Nov 2021 15:57
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 LabSampleId :
 SSTDCCC040

Quant Time: Nov 10 17:21:32 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 09 10:45:24 2021
 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	60	0.00
2	1,4-Dioxane	40.000	42.605	-6.5	63	0.00
3	Pyridine	40.000	41.724	-4.3	66	0.01
4	n-Nitrosodimethylamine	40.000	39.579	1.1	62	0.00
5 S	2-Fluorophenol	80.000	78.646	1.7	61	0.01
6	Aniline	40.000	38.912	2.7	59	0.00
7 S	Phenol-d6	80.000	77.697	2.9	59	0.00
8	2-Chlorophenol	40.000	39.857	0.4	61	0.01
9	Benzaldehyde	40.000	37.726	5.7	60	0.00
10 C	Phenol	40.000	38.933	2.7	59	0.00
11	bis(2-Chloroethyl)ether	40.000	37.755	5.6	57	0.00
12	1,3-Dichlorobenzene	40.000	39.342	1.6	61	0.01
13 C	1,4-Dichlorobenzene	40.000	39.192	2.0	61	0.00
14	1,2-Dichlorobenzene	40.000	39.890	0.3	61	0.00
15	Benzyl Alcohol	40.000	40.368	-0.9	59	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	36.646	8.4	56	0.01
17	2-Methylphenol	40.000	40.011	-0.0	60	0.00
18	Hexachloroethane	40.000	39.515	1.2	60	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	39.277	1.8	58	0.01
20	3+4-Methylphenols	40.000	39.508	1.2	58	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	58	0.00
22	Acetophenone	40.000	39.897	0.3	59	0.01
23 S	Nitrobenzene-d5	80.000	81.106	-1.4	59	0.01
24	Nitrobenzene	40.000	40.187	-0.5	59	0.01
25	Isophorone	40.000	39.446	1.4	57	0.00
26 C	2-Nitrophenol	40.000	41.540	-3.8	58	0.01
27	2,4-Dimethylphenol	40.000	39.546	1.1	57	0.00
28	bis(2-Chloroethoxy)methane	40.000	37.081	7.3	54	0.01
29 C	2,4-Dichlorophenol	40.000	40.900	-2.2	59	0.00
30	1,2,4-Trichlorobenzene	40.000	40.112	-0.3	59	0.00
31	Naphthalene	40.000	39.398	1.5	59	0.00
32	Benzoic acid	40.000	42.674	-6.7	56	0.00
33	4-Chloroaniline	40.000	39.559	1.1	58	0.00
34 C	Hexachlorobutadiene	40.000	41.792	-4.5	62	0.01
35	Caprolactam	40.000	41.385	-3.5	59	0.00
36 C	4-Chloro-3-methylphenol	40.000	40.892	-2.2	58	0.00
37	2-Methylnaphthalene	40.000	39.439	1.4	58	0.00
38	1-Methylnaphthalene	40.000	39.550	1.1	58	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	59	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	40.729	-1.8	61	0.01
41 P	Hexachlorocyclopentadiene	40.000	35.891	10.3	50	0.00
42 S	2,4,6-Tribromophenol	80.000	86.080	-7.6	64	0.00
43 C	2,4,6-Trichlorophenol	40.000	41.005	-2.5	59	0.00
44	2,4,5-Trichlorophenol	40.000	41.735	-4.3	61	0.00
45 S	2-Fluorobiphenyl	80.000	79.295	0.9	60	0.00
46	1,1'-Biphenyl	40.000	39.444	1.4	59	0.00
47	2-Chloronaphthalene	40.000	39.933	0.2	60	0.01

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111021\
 Data File : BP007921.D
 Acq On : 10 Nov 2021 15:57
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 LabSampleId :
 SSTDCCC040

Quant Time: Nov 10 17:21:32 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 09 10:45:24 2021
 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	41.988	-5.0	60	0.00
49	Acenaphthylene	40.000	40.091	-0.2	59	0.00
50	Dimethylphthalate	40.000	39.902	0.2	60	0.00
51	2,6-Dinitrotoluene	40.000	40.851	-2.1	59	0.00
52 C	Acenaphthene	40.000	39.625	0.9	59	0.00
53	3-Nitroaniline	40.000	40.898	-2.2	59	0.00
54 P	2,4-Dinitrophenol	40.000	43.530	-8.8	60	0.01
55	Dibenzofuran	40.000	37.364	6.6	56	0.00
56 P	4-Nitrophenol	40.000	40.243	-0.6	58	0.00
57	2,4-Dinitrotoluene	40.000	39.591	1.0	56	0.00
58	Fluorene	40.000	38.059	4.9	56	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	40.054	-0.1	58	0.00
60	Diethylphthalate	40.000	38.399	4.0	56	0.00
61	4-Chlorophenyl-phenylether	40.000	38.590	3.5	57	0.01
62	4-Nitroaniline	40.000	40.014	-0.0	57	0.00
63	Azobenzene	40.000	38.050	4.9	57	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	60	0.00
65	4,6-Dinitro-2-methylphenol	40.000	42.250	-5.6	57	0.00
66 c	n-Nitrosodiphenylamine	40.000	39.052	2.4	59	0.00
67	4-Bromophenyl-phenylether	40.000	40.862	-2.2	61	0.00
68	Hexachlorobenzene	40.000	40.938	-2.3	62	0.01
69	Atrazine	40.000	41.647	-4.1	62	0.00
70 C	Pentachlorophenol	40.000	42.996	-7.5	61	0.00
71	Phenanthrene	40.000	39.484	1.3	60	0.00
72	Anthracene	40.000	40.111	-0.3	61	0.00
73	Carbazole	40.000	39.905	0.2	61	0.00
74	Di-n-butylphthalate	40.000	40.182	-0.5	61	0.00
75 C	Fluoranthene	40.000	40.964	-2.4	63	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	67	0.00
77	Benzydine	40.000	36.982	7.5	55	0.00
78	Pyrene	40.000	37.685	5.8	63	0.00
79 S	Terphenyl-d14	80.000	77.086	3.6	66	0.00
80	Butylbenzylphthalate	40.000	38.999	2.5	65	0.00
81	Benzo(a)anthracene	40.000	39.029	2.4	67	0.00
82	3,3'-Dichlorobenzidine	40.000	39.339	1.7	66	0.00
83	Chrysene	40.000	39.135	2.2	67	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	38.419	4.0	65	0.00
85 c	Di-n-octyl phthalate	40.000	40.028	-0.1	66	0.00
86 I	Perylene-d12	20.000	20.000	0.0	67	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	38.184	4.5	65	0.01
88	Benzo(b)fluoranthene	40.000	41.169	-2.9	69	0.01
89	Benzo(k)fluoranthene	40.000	39.065	2.3	66	0.00
90 C	Benzo(a)pyrene	40.000	39.660	0.9	67	0.01
91	Dibenzo(a,h)anthracene	40.000	38.190	4.5	65	0.02
92	Benzo(g,h,i)perylene	40.000	37.816	5.5	64	0.01

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111021\
Data File : BP007921.D
Acq On : 10 Nov 2021 15:57
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Nov 10 17:21:32 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 09 10:45:24 2021
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