

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061121\
 Data File : BF124257.D
 Acq On : 11 Jun 2021 13:00
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jun 11 13:57:28 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 10 01:53:56 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	91	0.00
2	1,4-Dioxane	40.000	35.703	10.7	83	-0.02
3	Pyridine	40.000	36.109	9.7	84	-0.02
4	n-Nitrosodimethylamine	40.000	34.731	13.2	81	-0.02
5 S	2-Fluorophenol	80.000	75.803	5.2	88	-0.01
6	Aniline	40.000	37.580	6.1	90	0.00
7 S	Phenol-d6	80.000	73.244	8.4	87	0.00
8	2-Chlorophenol	40.000	39.894	0.3	92	-0.01
9	Benzaldehyde	40.000	38.402	4.0	94	0.00
10 C	Phenol	40.000	37.007	7.5	88	0.00
11	bis(2-Chloroethyl)ether	40.000	38.263	4.3	89	0.00
12	1,3-Dichlorobenzene	40.000	38.973	2.6	91	0.00
13 C	1,4-Dichlorobenzene	40.000	39.363	1.6	92	0.00
14	1,2-Dichlorobenzene	40.000	38.730	3.2	91	0.00
15	Benzyl Alcohol	40.000	38.053	4.9	91	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	35.323	11.7	85	0.00
17	2-Methylphenol	40.000	38.413	4.0	91	-0.01
18	Hexachloroethane	40.000	39.217	2.0	92	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	37.750	5.6	93	0.00
20	3+4-Methylphenols	40.000	39.093	2.3	92	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	99	0.00
22	Acetophenone	40.000	35.286	11.8	90	0.00
23 S	Nitrobenzene-d5	80.000	71.842	10.2	94	0.00
24	Nitrobenzene	40.000	35.518	11.2	90	0.00
25	Isophorone	40.000	35.075	12.3	92	0.00
26 C	2-Nitrophenol	40.000	35.061	12.3	90	0.00
27	2,4-Dimethylphenol	40.000	36.387	9.0	94	0.00
28	bis(2-Chloroethoxy)methane	40.000	34.149	14.6	92	0.00
29 C	2,4-Dichlorophenol	40.000	36.002	10.0	93	0.00
30	1,2,4-Trichlorobenzene	40.000	35.389	11.5	93	0.00
31	Naphthalene	40.000	35.428	11.4	92	0.00
32	Benzoic acid	40.000	35.031	12.4	96	0.00
33	4-Chloroaniline	40.000	36.835	7.9	96	-0.01
34 C	Hexachlorobutadiene	40.000	35.730	10.7	92	0.00
35	Caprolactam	40.000	36.003	10.0	95	0.00
36 C	4-Chloro-3-methylphenol	40.000	34.980	12.6	95	0.00
37	2-Methylnaphthalene	40.000	35.098	12.3	90	0.00
38	1-Methylnaphthalene	40.000	36.086	9.8	93	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	98	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	35.803	10.5	96	0.00
41 P	Hexachlorocyclopentadiene	40.000	34.113	14.7	90	-0.01
42 S	2,4,6-Tribromophenol	80.000	66.152	17.3	87	0.00
43 C	2,4,6-Trichlorophenol	40.000	34.935	12.7	90	0.00
44	2,4,5-Trichlorophenol	40.000	36.591	8.5	96	0.00
45 S	2-Fluorobiphenyl	80.000	72.781	9.0	95	0.00
46	1,1'-Biphenyl	40.000	35.590	11.0	95	0.00
47	2-Chloronaphthalene	40.000	35.419	11.5	95	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061121\
 Data File : BF124257.D
 Acq On : 11 Jun 2021 13:00
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 LabSampleId :
 SSTDCCC040

Quant Time: Jun 11 13:57:28 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 10 01:53:56 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	34.998	12.5	92	-0.01
49	Acenaphthylene	40.000	35.464	11.3	93	0.00
50	Dimethylphthalate	40.000	35.119	12.2	96	0.00
51	2,6-Dinitrotoluene	40.000	35.706	10.7	98	0.00
52 C	Acenaphthene	40.000	34.734	13.2	92	0.00
53	3-Nitroaniline	40.000	33.033	17.4	89	0.00
54 P	2,4-Dinitrophenol	40.000	36.759	8.1	100	0.00
55	Dibenzofuran	40.000	34.568	13.6	92	0.00
56 P	4-Nitrophenol	40.000	42.920	-7.3	108	0.00
57	2,4-Dinitrotoluene	40.000	34.666	13.3	88	0.00
58	Fluorene	40.000	35.644	10.9	93	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	34.302	14.2	93	0.00
60	Diethylphthalate	40.000	35.488	11.3	94	0.00
61	4-Chlorophenyl-phenylether	40.000	35.660	10.9	96	0.00
62	4-Nitroaniline	40.000	32.299	19.3	84	0.00
63	Azobenzene	40.000	34.024	14.9	92	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	92	0.00
65	4,6-Dinitro-2-methylphenol	40.000	34.204	14.5	84	0.00
66 c	n-Nitrosodiphenylamine	40.000	36.216	9.5	90	0.00
67	4-Bromophenyl-phenylether	40.000	35.088	12.3	91	0.00
68	Hexachlorobenzene	40.000	35.543	11.1	89	0.00
69	Atrazine	40.000	37.820	5.4	85	0.00
70 C	Pentachlorophenol	40.000	36.222	9.4	90	0.00
71	Phenanthrene	40.000	35.970	10.1	89	0.00
72	Anthracene	40.000	36.009	10.0	92	0.00
73	Carbazole	40.000	34.848	12.9	87	0.00
74	Di-n-butylphthalate	40.000	37.284	6.8	94	0.00
75 C	Fluoranthene	40.000	34.242	14.4	85	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	87	0.00
77	Benzydine	40.000	35.518	11.2	71	0.00
78	Pyrene	40.000	35.622	10.9	82	0.00
79 S	Terphenyl-d14	80.000	72.621	9.2	83	0.00
80	Butylbenzylphthalate	40.000	35.676	10.8	82	0.00
81	Benzo(a)anthracene	40.000	34.820	13.0	81	0.00
82	3,3'-Dichlorobenzidine	40.000	37.092	7.3	87	0.00
83	Chrysene	40.000	33.985	15.0	79	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	38.051	4.9	87	0.00
85 c	Di-n-octyl phthalate	40.000	42.012	-5.0	98	0.00
86 I	Perylene-d12	20.000	20.000	0.0	93	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	35.612	11.0	91	0.00
88	Benzo(b)fluoranthene	40.000	37.048	7.4	97	0.00
89	Benzo(k)fluoranthene	40.000	35.155	12.1	89	0.00
90 C	Benzo(a)pyrene	40.000	35.768	10.6	91	0.00
91	Dibenzo(a,h)anthracene	40.000	35.338	11.7	91	0.00
92	Benzo(g,h,i)perylene	40.000	35.027	12.4	88	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061121\
Data File : BF124257.D
Acq On : 11 Jun 2021 13:00
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_F
LabSampleId :
SSTDCCC040

Quant Time: Jun 11 13:57:28 2021
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF060321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 10 01:53:56 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