

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111020\
 Data File : BP003897.D
 Acq On : 10 Nov 2020 12:04
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 LabSampleId :
 SSTDCCC040

Quant Time: Nov 10 13:42:24 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 10 13:41:24 2020
 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	90	0.00
2	1,4-Dioxane	40.000	34.010	15.0	82	0.00
3	Pyridine	40.000	34.695	13.3	82	0.00
4	n-Nitrosodimethylamine	40.000	36.075	9.8	85	0.00
5 S	2-Fluorophenol	80.000	74.837	6.5	88	0.00
6	Aniline	40.000	36.569	8.6	85	0.00
7 S	Phenol-d6	80.000	73.457	8.2	86	0.00
8	2-Chlorophenol	40.000	37.639	5.9	88	0.00
9	Benzaldehyde	40.000	37.196	7.0	90	0.00
10 C	Phenol	40.000	36.439	8.9	86	0.00
11	bis(2-Chloroethyl)ether	40.000	35.957	10.1	85	0.00
12	1,3-Dichlorobenzene	40.000	37.136	7.2	89	0.00
13 C	1,4-Dichlorobenzene	40.000	37.092	7.3	88	0.00
14	1,2-Dichlorobenzene	40.000	37.257	6.9	88	0.00
15	Benzyl Alcohol	40.000	39.053	2.4	89	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	34.085	14.8	81	0.00
17	2-Methylphenol	40.000	37.404	6.5	87	0.00
18	Hexachloroethane	40.000	39.427	1.4	91	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	37.648	5.9	86	0.00
20	3+4-Methylphenols	40.000	37.704	5.7	87	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	87	0.00
22	Acetophenone	40.000	37.015	7.5	86	0.00
23 S	Nitrobenzene-d5	80.000	73.488	8.1	85	0.00
24	Nitrobenzene	40.000	37.464	6.3	87	0.00
25	Isophorone	40.000	38.118	4.7	87	0.00
26 C	2-Nitrophenol	40.000	44.203	-10.5	97	0.00
27	2,4-Dimethylphenol	40.000	38.143	4.6	89	0.00
28	bis(2-Chloroethoxy)methane	40.000	36.222	9.4	85	0.00
29 C	2,4-Dichlorophenol	40.000	39.857	0.4	91	0.00
30	1,2,4-Trichlorobenzene	40.000	38.171	4.6	90	0.00
31	Naphthalene	40.000	36.746	8.1	87	0.00
32	Benzoic acid	40.000	35.956	10.1	85	0.00
33	4-Chloroaniline	40.000	37.610	6.0	87	0.00
34 C	Hexachlorobutadiene	40.000	39.135	2.2	93	0.00
35	Caprolactam	40.000	41.695	-4.2	92	0.00
36 C	4-Chloro-3-methylphenol	40.000	38.683	3.3	89	0.00
37	2-Methylnaphthalene	40.000	37.087	7.3	87	0.00
38	1-Methylnaphthalene	40.000	36.724	8.2	87	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	88	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	37.629	5.9	89	0.00
41 P	Hexachlorocyclopentadiene	40.000	36.341	9.1	83	0.00
42 S	2,4,6-Tribromophenol	80.000	83.316	-4.1	95	0.00
43 C	2,4,6-Trichlorophenol	40.000	40.498	-1.2	93	0.00
44	2,4,5-Trichlorophenol	40.000	38.929	2.7	90	0.00
45 S	2-Fluorobiphenyl	80.000	74.186	7.3	89	0.00
46	1,1'-Biphenyl	40.000	36.652	8.4	88	0.00
47	2-Chloronaphthalene	40.000	36.523	8.7	87	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111020\
 Data File : BP003897.D
 Acq On : 10 Nov 2020 12:04
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_P
 LabSampleId :
 SSTDCCC040

Quant Time: Nov 10 13:42:24 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 10 13:41:24 2020
 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	40.711	-1.8	91	0.00
49	Acenaphthylene	40.000	37.132	7.2	88	0.00
50	Dimethylphthalate	40.000	37.231	6.9	88	0.00
51	2,6-Dinitrotoluene	40.000	39.172	2.1	90	0.00
52 C	Acenaphthene	40.000	37.300	6.8	89	0.00
53	3-Nitroaniline	40.000	39.329	1.7	91	0.00
54 P	2,4-Dinitrophenol	40.000	39.750	0.6	91	0.00
55	Dibenzofuran	40.000	36.558	8.6	88	0.00
56 P	4-Nitrophenol	40.000	39.968	0.1	89	0.00
57	2,4-Dinitrotoluene	40.000	40.711	-1.8	92	0.00
58	Fluorene	40.000	36.770	8.1	88	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	39.098	2.3	90	0.00
60	Diethylphthalate	40.000	37.853	5.4	90	0.00
61	4-Chlorophenyl-phenylether	40.000	37.387	6.5	90	0.00
62	4-Nitroaniline	40.000	40.592	-1.5	93	0.00
63	Azobenzene	40.000	35.863	10.3	84	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	89	0.00
65	4,6-Dinitro-2-methylphenol	40.000	38.198	4.5	93	0.00
66 c	n-Nitrosodiphenylamine	40.000	37.313	6.7	89	0.00
67	4-Bromophenyl-phenylether	40.000	38.505	3.7	92	0.00
68	Hexachlorobenzene	40.000	38.309	4.2	93	0.00
69	Atrazine	40.000	39.747	0.6	92	0.00
70 C	Pentachlorophenol	40.000	36.064	9.8	82	0.00
71	Phenanthrene	40.000	36.617	8.5	89	0.00
72	Anthracene	40.000	37.135	7.2	90	0.00
73	Carbazole	40.000	37.428	6.4	90	0.00
74	Di-n-butylphthalate	40.000	40.408	-1.0	93	0.00
75 C	Fluoranthene	40.000	37.181	7.0	89	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	86	0.00
77	Benzydine	40.000	45.199	-13.0	96	0.00
78	Pyrene	40.000	38.752	3.1	88	0.00
79 S	Terphenyl-d14	80.000	78.102	2.4	91	0.00
80	Butylbenzylphthalate	40.000	44.814	-12.0	96	0.00
81	Benzo(a)anthracene	40.000	37.582	6.0	87	0.00
82	3,3'-Dichlorobenzidine	40.000	40.688	-1.7	92	0.00
83	Chrysene	40.000	37.528	6.2	87	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	42.446	-6.1	93	0.00
85 c	Di-n-octyl phthalate	40.000	43.731	-9.3	95	0.00
86 I	Perylene-d12	20.000	20.000	0.0	90	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	36.009	10.0	90	0.00
88	Benzo(b)fluoranthene	40.000	36.410	9.0	93	0.00
89	Benzo(k)fluoranthene	40.000	33.777	15.6	83	0.00
90 C	Benzo(a)pyrene	40.000	35.796	10.5	89	0.00
91	Dibenzo(a,h)anthracene	40.000	36.077	9.8	89	0.00
92	Benzo(g,h,i)perylene	40.000	36.315	9.2	91	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111020\
Data File : BP003897.D
Acq On : 10 Nov 2020 12:04
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Nov 10 13:42:24 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\8270-BP110320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 10 13:41:24 2020
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