

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061121\
 Data File : BM030400.D
 Acq On : 11 Jun 2021 12:18
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC040

Quant Time: Jun 11 13:17:47 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 11 13:16:52 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	85	0.00
2	1,4-Dioxane	40.000	37.386	6.5	84	0.00
3	Pyridine	40.000	38.464	3.8	84	0.00
4	n-Nitrosodimethylamine	40.000	36.388	9.0	81	0.00
5 S	2-Fluorophenol	80.000	75.720	5.4	85	0.00
6	Aniline	40.000	36.059	9.9	81	0.00
7 S	Phenol-d6	80.000	74.745	6.6	82	0.00
8	2-Chlorophenol	40.000	36.826	7.9	84	0.00
9	Benzaldehyde	40.000	40.833	-2.1	84	0.00
10 C	Phenol	40.000	36.334	9.2	82	0.00
11	bis(2-Chloroethyl)ether	40.000	35.568	11.1	81	0.00
12	1,3-Dichlorobenzene	40.000	36.567	8.6	84	0.00
13 C	1,4-Dichlorobenzene	40.000	36.678	8.3	85	0.00
14	1,2-Dichlorobenzene	40.000	36.317	9.2	84	0.00
15	Benzyl Alcohol	40.000	35.741	10.6	79	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	33.940	15.2	76	0.00
17	2-Methylphenol	40.000	36.518	8.7	81	0.00
18	Hexachloroethane	40.000	36.344	9.1	82	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	34.675	13.3	76	0.00
20	3+4-Methylphenols	40.000	36.687	8.3	81	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	81	0.00
22	Acetophenone	40.000	37.375	6.6	79	0.00
23 S	Nitrobenzene-d5	80.000	76.144	4.8	79	0.00
24	Nitrobenzene	40.000	36.814	8.0	79	0.00
25	Isophorone	40.000	35.249	11.9	74	0.00
26 C	2-Nitrophenol	40.000	38.412	4.0	86	0.00
27	2,4-Dimethylphenol	40.000	36.480	8.8	78	0.00
28	bis(2-Chloroethoxy)methane	40.000	35.162	12.1	76	0.00
29 C	2,4-Dichlorophenol	40.000	38.237	4.4	82	0.00
30	1,2,4-Trichlorobenzene	40.000	37.228	6.9	82	0.00
31	Naphthalene	40.000	36.416	9.0	80	0.00
32	Benzoic acid	40.000	35.185	12.0	77	0.00
33	4-Chloroaniline	40.000	35.936	10.2	77	0.00
34 C	Hexachlorobutadiene	40.000	37.513	6.2	84	0.00
35	Caprolactam	40.000	33.072	17.3	68	0.00
36 C	4-Chloro-3-methylphenol	40.000	36.033	9.9	76	0.00
37	2-Methylnaphthalene	40.000	35.497	11.3	77	0.00
38	1-Methylnaphthalene	40.000	35.428	11.4	77	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	75	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	39.800	0.5	80	0.00
41 P	Hexachlorocyclopentadiene	40.000	39.048	2.4	79	0.00
42 S	2,4,6-Tribromophenol	80.000	74.778	6.5	72	0.00
43 C	2,4,6-Trichlorophenol	40.000	40.094	-0.2	80	0.00
44	2,4,5-Trichlorophenol	40.000	38.927	2.7	78	0.00
45 S	2-Fluorobiphenyl	80.000	76.807	4.0	77	0.00
46	1,1'-Biphenyl	40.000	38.159	4.6	76	0.00
47	2-Chloronaphthalene	40.000	37.367	6.6	76	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061121\
 Data File : BM030400.D
 Acq On : 11 Jun 2021 12:18
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC040

Quant Time: Jun 11 13:17:47 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 11 13:16:52 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	33.845	15.4	70	0.00
49	Acenaphthylene	40.000	37.129	7.2	74	0.00
50	Dimethylphthalate	40.000	35.228	11.9	71	0.00
51	2,6-Dinitrotoluene	40.000	36.734	8.2	72	0.00
52 C	Acenaphthene	40.000	35.850	10.4	73	0.00
53	3-Nitroaniline	40.000	33.104	17.2	69	0.00
54 P	2,4-Dinitrophenol	40.000	37.794	5.5	73	0.00
55	Dibenzofuran	40.000	35.803	10.5	74	0.00
56 P	4-Nitrophenol	40.000	35.517	11.2	68	0.00
57	2,4-Dinitrotoluene	40.000	34.441	13.9	71	0.00
58	Fluorene	40.000	35.460	11.3	71	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	37.318	6.7	74	0.00
60	Diethylphthalate	40.000	33.938	15.2	67	0.00
61	4-Chlorophenyl-phenylether	40.000	35.420	11.4	73	0.00
62	4-Nitroaniline	40.000	32.010	20.0	66	0.00
63	Azobenzene	40.000	34.042	14.9	66	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	69	0.00
65	4,6-Dinitro-2-methylphenol	40.000	40.195	-0.5	74	0.00
66 c	n-Nitrosodiphenylamine	40.000	37.705	5.7	70	0.00
67	4-Bromophenyl-phenylether	40.000	36.432	8.9	68	0.00
68	Hexachlorobenzene	40.000	37.906	5.2	72	0.00
69	Atrazine	40.000	43.297	-8.2	67	0.00
70 C	Pentachlorophenol	40.000	40.564	-1.4	71	0.00
71	Phenanthrene	40.000	36.552	8.6	68	0.00
72	Anthracene	40.000	37.164	7.1	68	0.00
73	Carbazole	40.000	36.331	9.2	66	0.00
74	Di-n-butylphthalate	40.000	36.857	7.9	62	0.00
75 C	Fluoranthene	40.000	36.223	9.4	67	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	69	0.00
77	Benzydine	40.000	48.557	-21.4	63	0.00
78	Pyrene	40.000	37.556	6.1	67	0.00
79 S	Terphenyl-d14	80.000	76.540	4.3	66	0.00
80	Butylbenzylphthalate	40.000	32.810	18.0	61	0.00
81	Benzo(a)anthracene	40.000	37.141	7.1	69	0.00
82	3,3'-Dichlorobenzidine	40.000	38.434	3.9	72	0.00
83	Chrysene	40.000	37.365	6.6	69	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	32.030	19.9	59	0.00
85 c	Di-n-octyl phthalate	40.000	33.906	15.2	64	0.00
86 I	Perylene-d12	20.000	20.000	0.0	81	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	40.221	-0.6	89	0.00
88	Benzo(b)fluoranthene	40.000	36.867	7.8	76	0.00
89	Benzo(k)fluoranthene	40.000	34.373	14.1	76	0.00
90 C	Benzo(a)pyrene	40.000	37.214	7.0	80	0.00
91	Dibenzo(a,h)anthracene	40.000	39.947	0.1	88	0.00
92	Benzo(g,h,i)perylene	40.000	41.173	-2.9	92	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061121\
Data File : BM030400.D
Acq On : 11 Jun 2021 12:18
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Jun 11 13:17:47 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060921.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 11 13:16:52 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