

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032921\
 Data File : BM029248.D
 Acq On : 29 Mar 2021 15:22
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC040

Quant Time: Mar 29 15:55:10 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM032621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 29 11:14:33 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	40.348	-0.9	94	0.00
3	Pyridine	40.000	43.643	-9.1	92	0.00
4	n-Nitrosodimethylamine	40.000	46.210	-15.5	95	0.00
5 S	2-Fluorophenol	80.000	81.705	-2.1	92	0.00
6	Aniline	40.000	42.255	-5.6	94	0.00
7 S	Phenol-d6	80.000	83.562	-4.5	92	0.00
8	2-Chlorophenol	40.000	41.002	-2.5	93	0.00
9	Benzaldehyde	40.000	40.069	-0.2	91	0.00
10 C	Phenol	40.000	42.007	-5.0	93	0.00
11	bis(2-Chloroethyl)ether	40.000	41.493	-3.7	91	0.00
12	1,3-Dichlorobenzene	40.000	40.313	-0.8	93	0.00
13 C	1,4-Dichlorobenzene	40.000	40.015	-0.0	91	0.00
14	1,2-Dichlorobenzene	40.000	39.889	0.3	91	0.00
15	Benzyl Alcohol	40.000	43.206	-8.0	93	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	43.140	-7.9	96	0.00
17	2-Methylphenol	40.000	41.901	-4.8	92	0.00
18	Hexachloroethane	40.000	41.321	-3.3	93	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	43.558	-8.9	92	0.00
20	3+4-Methylphenols	40.000	42.773	-6.9	92	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	90	0.00
22	Acetophenone	40.000	41.406	-3.5	92	0.00
23 S	Nitrobenzene-d5	80.000	84.416	-5.5	93	0.00
24	Nitrobenzene	40.000	42.162	-5.4	93	0.00
25	Isophorone	40.000	43.294	-8.2	93	0.00
26 C	2-Nitrophenol	40.000	43.479	-8.7	88	0.00
27	2,4-Dimethylphenol	40.000	41.490	-3.7	91	0.00
28	bis(2-Chloroethoxy)methane	40.000	42.744	-6.9	94	0.00
29 C	2,4-Dichlorophenol	40.000	41.949	-4.9	91	0.00
30	1,2,4-Trichlorobenzene	40.000	39.744	0.6	91	0.00
31	Naphthalene	40.000	40.572	-1.4	92	0.00
32	Benzoic acid	40.000	40.220	-0.5	88	0.00
33	4-Chloroaniline	40.000	41.807	-4.5	91	0.00
34 C	Hexachlorobutadiene	40.000	39.629	0.9	91	0.00
35	Caprolactam	40.000	39.816	0.5	88	0.00
36 C	4-Chloro-3-methylphenol	40.000	43.034	-7.6	91	0.00
37	2-Methylnaphthalene	40.000	41.271	-3.2	91	0.00
38	1-Methylnaphthalene	40.000	40.784	-2.0	91	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	88	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	38.769	3.1	88	0.00
41 P	Hexachlorocyclopentadiene	40.000	39.258	1.9	87	0.00
42 S	2,4,6-Tribromophenol	80.000	78.691	1.6	81	0.00
43 C	2,4,6-Trichlorophenol	40.000	41.215	-3.0	88	0.00
44	2,4,5-Trichlorophenol	40.000	41.425	-3.6	89	0.00
45 S	2-Fluorobiphenyl	80.000	80.977	-1.2	91	0.00
46	1,1'-Biphenyl	40.000	40.435	-1.1	90	0.00
47	2-Chloronaphthalene	40.000	39.828	0.4	89	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032921\
 Data File : BM029248.D
 Acq On : 29 Mar 2021 15:22
 Operator : CG/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC040

Quant Time: Mar 29 15:55:10 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM032621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 29 11:14:33 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	40.586	-1.5	91	0.00
49	Acenaphthylene	40.000	41.227	-3.1	90	0.00
50	Dimethylphthalate	40.000	40.983	-2.5	89	0.00
51	2,6-Dinitrotoluene	40.000	42.959	-7.4	91	0.00
52 C	Acenaphthene	40.000	39.924	0.2	88	0.00
53	3-Nitroaniline	40.000	43.517	-8.8	91	0.00
54 P	2,4-Dinitrophenol	40.000	37.710	5.7	84	0.00
55	Dibenzofuran	40.000	40.722	-1.8	90	0.00
56 P	4-Nitrophenol	40.000	41.400	-3.5	88	0.00
57	2,4-Dinitrotoluene	40.000	39.380	1.5	88	0.00
58	Fluorene	40.000	41.351	-3.4	90	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	41.248	-3.1	89	0.00
60	Diethylphthalate	40.000	42.383	-6.0	91	0.00
61	4-Chlorophenyl-phenylether	40.000	40.033	-0.1	88	0.00
62	4-Nitroaniline	40.000	40.545	-1.4	90	0.00
63	Azobenzene	40.000	44.655	-11.6	95	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	88	0.00
65	4,6-Dinitro-2-methylphenol	40.000	38.566	3.6	85	0.00
66 c	n-Nitrosodiphenylamine	40.000	40.968	-2.4	89	0.00
67	4-Bromophenyl-phenylether	40.000	36.663	8.3	80	0.00
68	Hexachlorobenzene	40.000	40.051	-0.1	88	0.00
69	Atrazine	40.000	42.734	-6.8	88	0.00
70 C	Pentachlorophenol	40.000	41.075	-2.7	87	0.00
71	Phenanthrene	40.000	40.808	-2.0	90	0.00
72	Anthracene	40.000	41.191	-3.0	89	0.00
73	Carbazole	40.000	41.573	-3.9	89	0.00
74	Di-n-butylphthalate	40.000	44.765	-11.9	91	0.00
75 C	Fluoranthene	40.000	41.440	-3.6	89	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	89	0.00
77	Benzydine	40.000	39.515	1.2	87	0.00
78	Pyrene	40.000	39.670	0.8	88	0.00
79 S	Terphenyl-d14	80.000	80.459	-0.6	88	0.00
80	Butylbenzylphthalate	40.000	40.501	-1.3	90	0.00
81	Benzo(a)anthracene	40.000	40.387	-1.0	89	0.00
82	3,3'-Dichlorobenzidine	40.000	41.502	-3.8	87	0.00
83	Chrysene	40.000	39.289	1.8	88	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	43.370	-8.4	96	0.00
85 c	Di-n-octyl phthalate	40.000	43.725	-9.3	97	0.00
86 I	Perylene-d12	20.000	20.000	0.0	89	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	40.972	-2.4	89	0.00
88	Benzo(b)fluoranthene	40.000	40.233	-0.6	87	0.00
89	Benzo(k)fluoranthene	40.000	41.075	-2.7	91	0.00
90 C	Benzo(a)pyrene	40.000	41.046	-2.6	89	0.00
91	Dibenzo(a,h)anthracene	40.000	41.221	-3.1	90	0.00
92	Benzo(g,h,i)perylene	40.000	40.611	-1.5	90	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM032921\
Data File : BM029248.D
Acq On : 29 Mar 2021 15:22
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Mar 29 15:55:10 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM032621.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Mar 29 11:14:33 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