

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110220\
 Data File : BM027668.D
 Acq On : 02 Nov 2020 09:19
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC040

Quant Time: Nov 02 10:16:21 2020
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM103020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 30 17:56:42 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	92	0.00
2	1,4-Dioxane	40.000	38.019	5.0	93	0.00
3	Pyridine	40.000	39.746	0.6	94	0.00
4	n-Nitrosodimethylamine	40.000	41.932	-4.8	103	0.00
5 S	2-Fluorophenol	80.000	77.263	3.4	94	0.00
6	Aniline	40.000	38.611	3.5	93	0.00
7 S	Phenol-d6	80.000	76.385	4.5	93	0.00
8	2-Chlorophenol	40.000	38.844	2.9	95	0.00
9	Benzaldehyde	40.000	40.406	-1.0	93	0.00
10 C	Phenol	40.000	38.449	3.9	95	0.00
11	bis(2-Chloroethyl)ether	40.000	38.243	4.4	94	0.00
12	1,3-Dichlorobenzene	40.000	38.879	2.8	96	0.00
13 C	1,4-Dichlorobenzene	40.000	38.289	4.3	94	0.00
14	1,2-Dichlorobenzene	40.000	38.609	3.5	96	0.00
15	Benzyl Alcohol	40.000	39.705	0.7	98	0.00
16	2,2'-oxybis(1-Chloropropane	40.000	38.214	4.5	95	0.00
17	2-Methylphenol	40.000	38.005	5.0	94	0.00
18	Hexachloroethane	40.000	39.355	1.6	98	0.00
19 P	n-Nitroso-di-n-propylamine	40.000	38.917	2.7	97	0.00
20	3+4-Methylphenols	40.000	39.598	1.0	97	0.00
21 I	Naphthalene-d8	20.000	20.000	0.0	96	0.00
22	Acetophenone	40.000	36.711	8.2	96	0.00
23 S	Nitrobenzene-d5	80.000	78.865	1.4	99	0.00
24	Nitrobenzene	40.000	38.794	3.0	100	0.00
25	Isophorone	40.000	37.081	7.3	96	0.00
26 C	2-Nitrophenol	40.000	40.460	-1.2	101	0.00
27	2,4-Dimethylphenol	40.000	36.974	7.6	95	0.00
28	bis(2-Chloroethoxy)methane	40.000	36.903	7.7	95	0.00
29 C	2,4-Dichlorophenol	40.000	38.117	4.7	98	0.00
30	1,2,4-Trichlorobenzene	40.000	38.120	4.7	100	0.00
31	Naphthalene	40.000	37.636	5.9	98	0.00
32	Benzoic acid	40.000	44.183	-10.5	118	0.01
33	4-Chloroaniline	40.000	36.963	7.6	95	0.00
34 C	Hexachlorobutadiene	40.000	37.299	6.8	97	0.00
35	Caprolactam	40.000	38.886	2.8	96	0.00
36 C	4-Chloro-3-methylphenol	40.000	38.657	3.4	97	0.00
37	2-Methylnaphthalene	40.000	37.666	5.8	99	0.00
38	1-Methylnaphthalene	40.000	37.943	5.1	98	0.00
39 I	Acenaphthene-d10	20.000	20.000	0.0	99	0.00
40	1,2,4,5-Tetrachlorobenzene	40.000	37.535	6.2	101	0.00
41 P	Hexachlorocyclopentadiene	40.000	38.773	3.1	103	0.00
42 S	2,4,6-Tribromophenol	80.000	78.143	2.3	102	0.00
43 C	2,4,6-Trichlorophenol	40.000	37.565	6.1	99	0.00
44	2,4,5-Trichlorophenol	40.000	36.918	7.7	98	0.00
45 S	2-Fluorobiphenyl	80.000	74.967	6.3	101	0.00
46	1,1'-Biphenyl	40.000	37.160	7.1	100	0.00
47	2-Chloronaphthalene	40.000	37.513	6.2	100	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110220\
 Data File : BM027668.D
 Acq On : 02 Nov 2020 09:19
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC040

Quant Time: Nov 02 10:16:21 2020
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM103020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 30 17:56:42 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	41.710	-4.3	105	0.00
49	Acenaphthylene	40.000	37.646	5.9	100	0.00
50	Dimethylphthalate	40.000	37.335	6.7	100	0.00
51	2,6-Dinitrotoluene	40.000	39.367	1.6	102	0.00
52 C	Acenaphthene	40.000	38.086	4.8	100	0.00
53	3-Nitroaniline	40.000	38.836	2.9	98	0.00
54 P	2,4-Dinitrophenol	40.000	44.185	-10.5	113	0.00
55	Dibenzofuran	40.000	37.379	6.6	101	0.00
56 P	4-Nitrophenol	40.000	37.425	6.4	98	0.00
57	2,4-Dinitrotoluene	40.000	39.710	0.7	102	0.00
58	Fluorene	40.000	37.800	5.5	100	0.00
59	2,3,4,6-Tetrachlorophenol	40.000	39.169	2.1	101	0.00
60	Diethylphthalate	40.000	37.745	5.6	101	0.00
61	4-Chlorophenyl-phenylether	40.000	37.462	6.3	103	0.00
62	4-Nitroaniline	40.000	40.030	-0.1	102	0.00
63	Azobenzene	40.000	38.243	4.4	102	0.00
64 I	Phenanthrene-d10	20.000	20.000	0.0	101	0.00
65	4,6-Dinitro-2-methylphenol	40.000	42.672	-6.7	112	0.00
66 c	n-Nitrosodiphenylamine	40.000	36.793	8.0	102	0.00
67	4-Bromophenyl-phenylether	40.000	36.771	8.1	102	0.00
68	Hexachlorobenzene	40.000	36.915	7.7	103	0.00
69	Atrazine	40.000	37.960	5.1	102	0.00
70 C	Pentachlorophenol	40.000	38.918	2.7	104	0.00
71	Phenanthrene	40.000	36.929	7.7	101	0.00
72	Anthracene	40.000	36.830	7.9	101	0.00
73	Carbazole	40.000	37.102	7.2	101	0.00
74	Di-n-butylphthalate	40.000	37.858	5.4	104	0.00
75 C	Fluoranthene	40.000	37.115	7.2	102	0.00
76 I	Chrysene-d12	20.000	20.000	0.0	103	0.00
77	Benzydine	40.000	36.669	8.3	98	0.00
78	Pyrene	40.000	37.235	6.9	103	0.00
79 S	Terphenyl-d14	80.000	74.158	7.3	104	0.00
80	Butylbenzylphthalate	40.000	39.772	0.6	107	0.00
81	Benzo(a)anthracene	40.000	37.429	6.4	104	0.00
82	3,3'-Dichlorobenzidine	40.000	37.300	6.8	101	0.00
83	Chrysene	40.000	37.145	7.1	104	0.00
84	Bis(2-ethylhexyl)phthalate	40.000	37.959	5.1	105	0.00
85 c	Di-n-octyl phthalate	40.000	38.598	3.5	107	0.00
86 I	Perylene-d12	20.000	20.000	0.0	105	0.00
87	Indeno(1,2,3-cd)pyrene	40.000	36.088	9.8	104	0.00
88	Benzo(b)fluoranthene	40.000	35.123	12.2	103	0.00
89	Benzo(k)fluoranthene	40.000	36.834	7.9	105	0.00
90 C	Benzo(a)pyrene	40.000	35.764	10.6	103	0.00
91	Dibenzo(a,h)anthracene	40.000	35.447	11.4	103	0.00
92	Benzo(g,h,i)perylene	40.000	35.572	11.1	102	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110220\
Data File : BM027668.D
Acq On : 02 Nov 2020 09:19
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Nov 02 10:16:21 2020
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM103020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Oct 30 17:56:42 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