

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082120\
 Data File : BM027174.D
 Acq On : 21 Aug 2020 17:50
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC040

Quant Time: Aug 24 10:42:45 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082120.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 24 10:33:46 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	84	0.01
2	1,4-Dioxane	0.490	0.390	20.4	73	0.00
3	Pyridine	1.376	1.164	15.4	74	0.00
4	n-Nitrosodimethylamine	0.600	0.553	7.8	83	0.00
5 S	2-Fluorophenol	1.026	0.948	7.6	80	0.00
6	Aniline	1.600	1.529	4.4	113	0.01
7 S	Phenol-d6	1.409	1.346	4.5	111	0.00
8	2-Chlorophenol	1.068	1.055	1.2	115	0.01
9	Benzaldehyde	0.964	0.877	9.0	109	0.01
10 C	Phenol	1.342	1.270	5.4	112	0.00
11	bis(2-Chloroethyl)ether	1.097	1.061	3.3	115	0.00
12	1,3-Dichlorobenzene	1.494	1.400	6.3	87	0.00
13 C	1,4-Dichlorobenzene	1.508	1.395	7.5	85	0.01
14	1,2-Dichlorobenzene	1.430	1.348	5.7	87	0.01
15	Benzyl Alcohol	1.280	1.277	0.2	89	0.00
16	2,2'-oxybis(1-Chloropropane	0.880	0.828	5.9	84	0.00
17	2-Methylphenol	0.929	0.910	2.0	88	0.00
18	Hexachloroethane	0.612	0.621	-1.5	94	0.01
19 P	n-Nitroso-di-n-propylamine	1.081	1.116	-3.2	92	0.01
20	3+4-Methylphenols	1.243	1.243	0.0	88	0.00
21 I	Naphthalene-d8	1.000	1.000	0.0	93	0.01
22	Acetophenone	0.538	0.492	8.6	93	0.01
23 S	Nitrobenzene-d5	0.485	0.441	9.1	91	0.00
24	Nitrobenzene	0.491	0.446	9.2	92	0.00
25	Isophorone	0.710	0.658	7.3	91	0.00
26 C	2-Nitrophenol	0.153	0.149	2.6	96	0.00
27	2,4-Dimethylphenol	0.238	0.221	7.1	92	0.00
28	bis(2-Chloroethoxy)methane	0.433	0.418	3.5	95	0.00
29 C	2,4-Dichlorophenol	0.332	0.329	0.9	98	0.00
30	1,2,4-Trichlorobenzene	0.435	0.410	5.7	97	0.00
31	Naphthalene	1.014	0.954	5.9	95	0.00
32	Benzoic acid	0.170	0.177	-4.1	93	0.00
33	4-Chloroaniline	0.419	0.424	-1.2	98	0.01
34 C	Hexachlorobutadiene	0.314	0.304	3.2	103	0.01
35	Caprolactam	0.090	0.097	-7.8	100	0.00
36 C	4-Chloro-3-methylphenol	0.343	0.361	-5.2	100	0.00
37	2-Methylnaphthalene	0.768	0.762	0.8	99	0.01
38	1-Methylnaphthalene	0.739	0.739	0.0	100	0.00
39 I	Acenaphthene-d10	1.000	1.000	0.0	107	0.00
40	1,2,4,5-Tetrachlorobenzene	0.704	0.604	14.2	105	0.01
41 P	Hexachlorocyclopentadiene	0.389	0.351	9.8	106	0.01
42 S	2,4,6-Tribromophenol	0.324	0.335	-3.4	118	0.00
43 C	2,4,6-Trichlorophenol	0.435	0.395	9.2	106	0.00
44	2,4,5-Trichlorophenol	0.461	0.416	9.8	107	0.00

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082120\
 Data File : BM027174.D
 Acq On : 21 Aug 2020 17:50
 Operator : JU/CG
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 LabSampleId :
 SSTDCCC040

Quant Time: Aug 24 10:42:45 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082120.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 24 10:33:46 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	AvqRF	CCRF	%Dev	Area%	Dev(min)
45 S	2-Fluorobiphenyl	1.523	1.337	12.2	106	0.00
46	1,1'-Biphenyl	1.499	1.322	11.8	105	0.00
47	2-Chloronaphthalene	1.278	1.129	11.7	106	0.00
48	2-Nitroaniline	0.408	0.391	4.2	106	0.00
49	Acenaphthylene	1.777	1.665	6.3	107	0.00
50	Dimethylphthalate	1.624	1.511	7.0	108	0.00
51	2,6-Dinitrotoluene	0.323	0.311	3.7	107	0.00
52 C	Acenaphthene	1.128	1.062	5.9	107	0.00
53	3-Nitroaniline	0.311	0.314	-1.0	105	0.00
54 P	2,4-Dinitrophenol	0.169	0.175	-3.6	115	0.00
55	Dibenzofuran	1.880	1.800	4.3	110	0.00
56 P	4-Nitrophenol	0.248	0.258	-4.0	112	0.00
57	2,4-Dinitrotoluene	0.449	0.457	-1.8	110	0.00
58	Fluorene	1.552	1.536	1.0	112	0.00
59	2,3,4,6-Tetrachlorophenol	0.452	0.440	2.7	111	0.00
60	Diethylphthalate	1.611	1.614	-0.2	112	0.00
61	4-Chlorophenyl-phenylether	0.891	0.860	3.5	113	0.00
62	4-Nitroaniline	0.349	0.374	-7.2	112	0.00
63	Azobenzene	1.653	1.681	-1.7	113	0.01
64 I	Phenanthrene-d10	1.000	1.000	0.0	115	0.01
65	4,6-Dinitro-2-methylphenol	0.098	0.092	6.1	115	0.00
66 c	n-Nitrosodiphenylamine	0.568	0.517	9.0	113	0.00
67	4-Bromophenyl-phenylether	0.241	0.222	7.9	117	0.00
68	Hexachlorobenzene	0.287	0.263	8.4	117	0.00
69	Atrazine	0.215	0.204	5.1	117	0.00
70 C	Pentachlorophenol	0.155	0.153	1.3	120	0.00
71	Phenanthrene	1.112	1.046	5.9	116	0.00
72	Anthracene	1.063	1.021	4.0	117	0.00
73	Carbazole	0.978	0.949	3.0	118	0.00
74	Di-n-butylphthalate	1.129	1.178	-4.3	124	0.00
75 C	Fluoranthene	1.389	1.346	3.1	120	0.01
76 I	Chrysene-d12	1.000	1.000	0.0	129	0.00
77	Benzidine	0.347	0.440	-26.8#	150	0.00
78	Pyrene	1.217	1.110	8.8	122	0.00
79 S	Terphenyl-d14	1.071	0.987	7.8	125	0.00
80	Butylbenzylphthalate	0.445	0.456	-2.5	130	0.00
81	Benzo(a)anthracene	1.296	1.219	5.9	128	0.00
82	3,3'-Dichlorobenzidine	0.452	0.450	0.4	131	0.00
83	Chrysene	1.259	1.172	6.9	128	0.00
84	Bis(2-ethylhexyl)phthalate	0.633	0.665	-5.1	132	0.00
85 c	Di-n-octyl phthalate	1.044	1.135	-8.7	137	0.00
86 I	Perylene-d12	1.000	1.000	0.0	134	0.00
87	Indeno(1,2,3-cd)pyrene	1.565	1.397	10.7	133	0.02

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM082120\
Data File : BM027174.D
Acq On : 21 Aug 2020 17:50
Operator : JU/CG
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Aug 24 10:42:45 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM082120.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 24 10:33:46 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.334	1.282	3.9	132	0.00
89	Benzo(k)fluoranthene	1.306	1.184	9.3	129	0.00
90 C	Benzo(a)pyrene	1.218	1.145	6.0	132	0.00
91	Dibenzo(a,h)anthracene	1.303	1.163	10.7	134	0.02
92	Benzo(g,h,i)perylene	1.255	1.115	11.2	134	0.02

(#) = Out of Range

SPCC's out = 0 CCC's out = 0