

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF041223\
 Data File : BF132868.D
 Acq On : 12 Apr 2023 19:21
 Operator : CG\JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Quant Time: Apr 13 05:21:06 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF041223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 13 05:14:42 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.769	152	250486	20.000	ng	0.00
21) Naphthalene-d8	8.057	136	924097	20.000	ng	0.00
39) Acenaphthene-d10	9.810	164	451177	20.000	ng	0.00
64) Phenanthrene-d10	11.286	188	792492	20.000	ng	0.00
76) Chrysene-d12	13.922	240	514295	20.000	ng	0.00
86) Perylene-d12	15.345	264	407320	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	1789089	109.533	ng	0.00
7) Phenol-d6	6.404	99	2278468	109.349	ng	0.00
23) Nitrobenzene-d5	7.345	82	2009207	114.659	ng	0.00
42) 2,4,6-Tribromophenol	10.598	330	533436	124.702	ng	0.00
45) 2-Fluorobiphenyl	9.139	172	3069228	103.857	ng	0.00
79) Terphenyl-d14	12.880	244	3483397	118.142	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.463	88	459316	56.571	ng	99
3) Pyridine	3.181	79	1280099	57.705	ng	98
4) n-Nitrosodimethylamine	3.146	42	606458	58.069	ng	97
6) Aniline	6.440	93	1540046	57.303	ng	100
8) 2-Chlorophenol	6.557	128	892889	54.754	ng	96
9) Benzaldehyde	6.316	77	613023	58.829	ng	97
10) Phenol	6.422	94	1282087	54.406	ng	92
11) bis(2-Chloroethyl)ether	6.516	93	948515	55.227	ng	97
12) 1,3-Dichlorobenzene	6.716	146	968078	54.608	ng	98
13) 1,4-Dichlorobenzene	6.793	146	966020	54.188	ng	97
14) 1,2-Dichlorobenzene	6.945	146	877067	52.705	ng	98
15) Benzyl Alcohol	6.916	79	797066	58.507	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.057	45	1575986	55.484	ng	99
17) 2-Methylphenol	7.022	107	826754	56.291	ng	99
18) Hexachloroethane	7.287	117	336572	55.433	ng	99
19) n-Nitroso-di-n-propyla...	7.198	70	682157	56.987	ng	99
20) 3+4-Methylphenols	7.187	107	923412	52.234	ng	94
22) Acetophenone	7.187	105	1165277	53.585	ng	# 88
24) Nitrobenzene	7.363	77	1025773	57.454	ng	99
25) Isophorone	7.598	82	1883234	59.246	ng	97
26) 2-Nitrophenol	7.675	139	429008	68.121	ng	97
27) 2,4-Dimethylphenol	7.716	122	778514	57.795	ng	100
28) bis(2-Chloroethoxy)met...	7.810	93	1090601	57.358	ng	100
29) 2,4-Dichlorophenol	7.916	162	732953	58.686	ng	99
30) 1,2,4-Trichlorobenzene	7.998	180	762612	56.340	ng	99
31) Naphthalene	8.081	128	2551544	54.707	ng	99
32) Benzoic acid	7.851	122	428280m	60.581	ng	
33) 4-Chloroaniline	8.128	127	1089186	57.603	ng	96
34) Hexachlorobutadiene	8.198	225	429489	57.418	ng	98
35) Caprolactam	8.516	113	246965	67.630	ng	98
36) 4-Chloro-3-methylphenol	8.610	107	788354	59.711	ng	98
37) 2-Methylnaphthalene	8.769	142	1704286	54.341	ng	98
38) 1-Methylnaphthalene	8.869	142	1580645	53.973	ng	99
40) 1,2,4,5-Tetrachloroben...	8.939	216	718543	56.424	ng	100
41) Hexachlorocyclopentadiene	8.928	237	393272	64.440	ng	100
43) 2,4,6-Trichlorophenol	9.045	196	473105	60.458	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF041223\
 Data File : BF132868.D
 Acq On : 12 Apr 2023 19:21
 Operator : CG\JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Quant Time: Apr 13 05:21:06 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF041223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 13 05:14:42 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.081	196	550425	59.383	ng	96
46) 1,1'-Biphenyl	9.234	154	1976525	54.265	ng	100
47) 2-Chloronaphthalene	9.257	162	1462239	55.133	ng	96
48) 2-Nitroaniline	9.351	65	499852	60.475	ng	97
49) Acenaphthylene	9.675	152	2347161	54.428	ng	100
50) Dimethylphthalate	9.539	163	1723467	57.619	ng	99
51) 2,6-Dinitrotoluene	9.598	165	408964	62.308	ng	91
52) Acenaphthene	9.845	154	1540587	55.611	ng	100
53) 3-Nitroaniline	9.763	138	494523	63.843	ng	98
54) 2,4-Dinitrophenol	9.863	184	184107	61.068	ng	91
55) Dibenzofuran	10.016	168	2118077	54.017	ng	97
56) 4-Nitrophenol	9.916	139	381060	62.807	ng	89
57) 2,4-Dinitrotoluene	9.998	165	525404	64.557	ng	99
58) Fluorene	10.357	166	1553259	53.086	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.134	232	432430	60.676	ng	99
60) Diethylphthalate	10.239	149	1610212	59.762	ng	99
61) 4-Chlorophenyl-phenyle...	10.351	204	755323	54.160	ng	97
62) 4-Nitroaniline	10.381	138	485976	64.117	ng	96
63) Azobenzene	10.510	77	1807127	56.098	ng	98
65) 4,6-Dinitro-2-methylph...	10.404	198	243411	61.252	ng	98
66) n-Nitrosodiphenylamine	10.475	169	1433661	56.835	ng	100
67) 4-Bromophenyl-phenylether	10.839	248	498722	58.807	ng	94
68) Hexachlorobenzene	10.904	284	520229	58.727	ng	94
69) Atrazine	10.998	200	412250	62.619	ng	99
70) Pentachlorophenol	11.092	266	374170	64.530	ng	99
71) Phenanthrene	11.316	178	2409329	55.040	ng	98
72) Anthracene	11.369	178	2462839	55.066	ng	99
73) Carbazole	11.522	167	2201305	55.859	ng	99
74) Di-n-butylphthalate	11.857	149	2223605	67.370	ng	99
75) Fluoranthene	12.498	202	2490865	56.799	ng	99
77) Benzidine	12.622	184	456040	62.136	ng	100
78) Pyrene	12.727	202	2567254	59.800	ng	98
80) Butylbenzylphthalate	13.351	149	466757	62.602	ng	99
81) Benzo(a)anthracene	13.910	228	2078749	58.701	ng	99
82) 3,3'-Dichlorobenzidine	13.880	252	492926	61.484	ng	99
83) Chrysene	13.951	228	2017488	58.427	ng	97
84) Bis(2-ethylhexyl)phtha...	13.916	149	630336	63.217	ng	99
85) Di-n-octyl phthalate	14.527	149	964068	63.366	ng	96
87) Indeno(1,2,3-cd)pyrene	16.751	276	1657397	60.487	ng	100
88) Benzo(b)fluoranthene	14.939	252	1641240	63.599	ng	100
89) Benzo(k)fluoranthene	14.974	252	1505766	56.902	ng	98
90) Benzo(a)pyrene	15.292	252	1447194	61.436	ng	99
91) Dibenzo(a,h)anthracene	16.774	278	1363535	60.811	ng	98
92) Benzo(g,h,i)perylene	17.174	276	1418110	60.835	ng	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

SSTDICC060

Response via : Initial Calibration

