

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102424\
 Data File : BF139997.D
 Acq On : 24 Oct 2024 13:21
 Operator : RC/JU
 Sample : PB164208BS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB164208BS

Quant Time: Oct 24 13:41:52 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101824.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 15:07:50 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.892	152	146768	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	559393	20.000	ng	0.00
39) Acenaphthene-d10	9.927	164	307081	20.000	ng	0.00
64) Phenanthrene-d10	11.416	188	545635	20.000	ng	0.00
76) Chrysene-d12	14.057	240	272685	20.000	ng	0.00
86) Perylene-d12	15.527	264	290439	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.516	112	1173430	125.163	ng	0.02
7) Phenol-d6	6.516	99	1488768	122.609	ng	0.00
23) Nitrobenzene-d5	7.451	82	961002	95.214	ng	0.00
42) 2,4,6-Tribromophenol	10.722	330	445211	154.999	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	1726013	92.893	ng	0.00
79) Terphenyl-d14	12.998	244	1757783	105.040	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.763	88	164401	37.610	ng	94
3) Pyridine	3.522	79	434284	40.082	ng	94
4) n-Nitrosodimethylamine	3.469	42	237180	40.744	ng	94
6) Aniline	6.551	93	457846	40.458	ng	98
8) 2-Chlorophenol	6.675	128	456067	47.585	ng	97
9) Benzaldehyde	6.439	77	109292	16.000	ng	98
10) Phenol	6.528	94	560448m	44.770	ng	
11) bis(2-Chloroethyl)ether	6.628	93	441357	45.615	ng	98
12) 1,3-Dichlorobenzene	6.834	146	487132	44.277	ng	98
13) 1,4-Dichlorobenzene	6.910	146	492253	44.784	ng	98
14) 1,2-Dichlorobenzene	7.057	146	477151	46.512	ng	99
15) Benzyl Alcohol	7.028	79	411790	46.233	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.163	45	714292	43.295	ng	97
17) 2-Methylphenol	7.134	107	374494	45.823	ng	100
18) Hexachloroethane	7.404	117	175746	44.837	ng	97
19) n-Nitroso-di-n-propyla...	7.304	70	329758	44.778	ng	96
20) 3+4-Methylphenols	7.286	107	460785	44.080	ng	94
22) Acetophenone	7.298	105	614828	44.873	ng	98
24) Nitrobenzene	7.469	77	496780	44.893	ng	100
25) Isophorone	7.710	82	894254	46.681	ng	99
26) 2-Nitrophenol	7.786	139	238511	57.133	ng	95
27) 2,4-Dimethylphenol	7.816	122	378965	54.441	ng	99
28) bis(2-Chloroethoxy)met...	7.916	93	533558	45.971	ng	99
29) 2,4-Dichlorophenol	8.022	162	377651	47.630	ng	99
30) 1,2,4-Trichlorobenzene	8.110	180	399159	45.497	ng	99
31) Naphthalene	8.192	128	1311180	45.448	ng	100
32) Benzoic acid	7.933	122	293183	48.289	ng	95
33) 4-Chloroaniline	8.239	127	217150	22.074	ng	97
34) Hexachlorobutadiene	8.310	225	260021	47.170	ng	98
35) Caprolactam	8.610	113	120588m	47.832	ng	
36) 4-Chloro-3-methylphenol	8.716	107	407211	46.382	ng	97
37) 2-Methylnaphthalene	8.886	142	828427	46.886	ng	100
38) 1-Methylnaphthalene	8.986	142	770255	44.433	ng	100
40) 1,2,4,5-Tetrachloroben...	9.051	216	399655	48.241	ng	99
41) Hexachlorocyclopentadiene	9.039	237	516673	179.237	ng	99
43) 2,4,6-Trichlorophenol	9.163	196	296773	50.597	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.198	196	289527	47.844	ng	99
46) 1,1'-Biphenyl	9.351	154	1005890	47.054	ng	99
47) 2-Chloronaphthalene	9.375	162	800201	46.476	ng	99
48) 2-Nitroaniline	9.469	65	270414	51.352	ng	91
49) Acenaphthylene	9.792	152	1236148	49.661	ng	100
50) Dimethylphthalate	9.651	163	921402	48.172	ng	100
51) 2,6-Dinitrotoluene	9.710	165	204338	49.339	ng	96
52) Acenaphthene	9.963	154	870187	54.042	ng	98
53) 3-Nitroaniline	9.875	138	134327	31.452	ng	97
54) 2,4-Dinitrophenol	9.986	184	225831	127.062	ng	# 1
55) Dibenzofuran	10.133	168	1088550	47.118	ng	99
56) 4-Nitrophenol	10.033	139	343997	104.692	ng	97
57) 2,4-Dinitrotoluene	10.116	165	275731	54.140	ng	95
58) Fluorene	10.480	166	823797	46.816	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.251	232	246162	51.890	ng	97
60) Diethylphthalate	10.351	149	892502	47.523	ng	99
61) 4-Chlorophenyl-phenyle...	10.469	204	421137	47.392	ng	98
62) 4-Nitroaniline	10.492	138	200030	49.590	ng	95
63) Azobenzene	10.627	77	917547	46.084	ng	98
65) 4,6-Dinitro-2-methylph...	10.522	198	151935	68.266	ng	91
66) n-Nitrosodiphenylamine	10.586	169	769526	47.121	ng	100
67) 4-Bromophenyl-phenylether	10.963	248	275211	48.961	ng	93
68) Hexachlorobenzene	11.027	284	307480	48.638	ng	94
69) Atrazine	11.116	200	251414	56.833	ng	99
70) Pentachlorophenol	11.216	266	383001	99.824	ng	100
71) Phenanthrene	11.439	178	1222910	47.448	ng	100
72) Anthracene	11.492	178	1239887	49.306	ng	100
73) Carbazole	11.645	167	1069883	45.747	ng	99
74) Di-n-butylphthalate	11.974	149	1293914	47.887	ng	100
75) Fluoranthene	12.627	202	1212798	46.548	ng	100
77) Benzidine	12.745	184	229171	51.110	ng	99
78) Pyrene	12.857	202	1214033	50.862	ng	100
80) Butylbenzylphthalate	13.474	149	390860	54.620	ng	99
81) Benzo(a)anthracene	14.045	228	885407	49.880	ng	100
82) 3,3'-Dichlorobenzidine	14.004	252	193751	37.436	ng	99
83) Chrysene	14.080	228	793014	48.725	ng	99
84) Bis(2-ethylhexyl)phtha...	14.033	149	433847	54.037	ng	# 99
85) Di-n-octyl phthalate	14.645	149	716965	49.097	ng	96
87) Indeno(1,2,3-cd)pyrene	17.033	276	1042017	55.769	ng	98
88) Benzo(b)fluoranthene	15.098	252	764657	43.220	ng	99
89) Benzo(k)fluoranthene	15.127	252	775417	50.820	ng	100
90) Benzo(a)pyrene	15.468	252	763391	52.439	ng	99
91) Dibenzo(a,h)anthracene	17.051	278	854677	54.832	ng	98
92) Benzo(g,h,i)perylene	17.486	276	797561	51.226	ng	98

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

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