

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010325\
 Data File : BF141065.D
 Acq On : 03 Jan 2025 14:03
 Operator : RC/JU
 Sample : PB165945BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB165945BS

Quant Time: Jan 03 14:24:12 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 27 00:31:16 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.828	152	261678	20.000	ng	0.00
21) Naphthalene-d8	8.110	136	1044303	20.000	ng	0.00
39) Acenaphthene-d10	9.863	164	561214	20.000	ng	0.00
64) Phenanthrene-d10	11.345	188	930135	20.000	ng	0.00
76) Chrysene-d12	13.992	240	595107	20.000	ng	0.00
86) Perylene-d12	15.457	264	551132	20.000	ng	0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.451	112	2192753	126.312	ng	0.02
7) Phenol-d6	6.457	99	2777986	124.221	ng	0.00
23) Nitrobenzene-d5	7.392	82	1773872	108.674	ng	0.00
42) 2,4,6-Tribromophenol	10.651	330	859584	157.625	ng	0.00
45) 2-Fluorobiphenyl	9.186	172	3090607	87.740	ng	0.00
79) Terphenyl-d14	12.933	244	3453233	96.383	ng	0.00

Target Compounds					Qvalue
2) 1,4-Dioxane	2.663	88	308273	39.224	ng 98
3) Pyridine	3.399	79	813345	42.346	ng 99
4) n-Nitrosodimethylamine	3.352	42	434558	44.654	ng 99
6) Aniline	6.487	93	971953	48.732	ng 95
8) 2-Chlorophenol	6.610	128	904395	49.413	ng 97
9) Benzaldehyde	6.375	77	156674	7.635	ng 97
10) Phenol	6.469	94	1104713	47.773	ng 94
11) bis(2-Chloroethyl)ether	6.563	93	850232	45.537	ng 98
12) 1,3-Dichlorobenzene	6.769	146	918544	45.475	ng 99
13) 1,4-Dichlorobenzene	6.845	146	928083	45.578	ng 99
14) 1,2-Dichlorobenzene	6.998	146	885030	46.617	ng 98
15) Benzyl Alcohol	6.969	79	770181	47.926	ng 99
16) 2,2'-oxybis(1-Chloropr...	7.104	45	1297222	45.941	ng 98
17) 2-Methylphenol	7.075	107	726082	47.752	ng 100
18) Hexachloroethane	7.340	117	338952	46.859	ng 100
19) n-Nitroso-di-n-propyla...	7.245	70	621530	45.366	ng 97
20) 3+4-Methylphenols	7.234	107	855069	44.438	ng # 80
22) Acetophenone	7.239	105	1147179	44.780	ng 98
24) Nitrobenzene	7.410	77	924281	50.676	ng 96
25) Isophorone	7.651	82	1687984	47.523	ng 99
26) 2-Nitrophenol	7.728	139	447221	58.525	ng 97
27) 2,4-Dimethylphenol	7.763	122	728497	56.119	ng 99
28) bis(2-Chloroethoxy)met...	7.863	93	1031419	45.832	ng 100
29) 2,4-Dichlorophenol	7.963	162	719786	48.890	ng 100
30) 1,2,4-Trichlorobenzene	8.051	180	755576	45.342	ng 99
31) Naphthalene	8.134	128	2500894	45.447	ng 100
32) Benzoic acid	7.881	122	554184	51.255	ng 96
33) 4-Chloroaniline	8.175	127	493711	24.800	ng 99
34) Hexachlorobutadiene	8.251	225	455803	46.398	ng 99
35) Caprolactam	8.545	113	244555m	48.861	ng
36) 4-Chloro-3-methylphenol	8.657	107	780837	47.757	ng 98
37) 2-Methylnaphthalene	8.822	142	1626476	46.202	ng 100
38) 1-Methylnaphthalene	8.922	142	1522459	44.005	ng 99
40) 1,2,4,5-Tetrachloroben...	8.986	216	723909	46.886	ng 99
41) Hexachlorocyclopentadiene	8.975	237	909341	177.726	ng 100
43) 2,4,6-Trichlorophenol	9.098	196	530051	51.582	ng 99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.133	196	527631	50.183	ng	99
46) 1,1'-Biphenyl	9.286	154	1978272	46.036	ng	99
47) 2-Chloronaphthalene	9.310	162	1501403	45.050	ng	100
48) 2-Nitroaniline	9.404	65	484004	54.786	ng	92
49) Acenaphthylene	9.728	152	2345475	48.888	ng	99
50) Dimethylphthalate	9.586	163	1792105	48.547	ng	100
51) 2,6-Dinitrotoluene	9.645	165	408682	53.808	ng	95
52) Acenaphthene	9.898	154	1691822	52.161	ng	100
53) 3-Nitroaniline	9.810	138	293424	36.667	ng	# 93
54) 2,4-Dinitrophenol	9.922	184	403862	113.034	ng	# 1
55) Dibenzofuran	10.069	168	2127512	46.675	ng	100
56) 4-Nitrophenol	9.975	139	701887	111.154	ng	94
57) 2,4-Dinitrotoluene	10.051	165	542028	57.822	ng	95
58) Fluorene	10.410	166	1593297	45.077	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.186	232	465819	53.494	ng	98
60) Diethylphthalate	10.286	149	1726921	47.386	ng	100
61) 4-Chlorophenyl-phenyle...	10.404	204	784094	46.102	ng	98
62) 4-Nitroaniline	10.428	138	430938	52.362	ng	95
63) Azobenzene	10.563	77	1750610	45.765	ng	99
65) 4,6-Dinitro-2-methylph...	10.457	198	279427	70.024	ng	94
66) n-Nitrosodiphenylamine	10.522	169	1481992	50.598	ng	99
67) 4-Bromophenyl-phenylether	10.892	248	523658	52.080	ng	99
68) Hexachlorobenzene	10.957	284	574476	51.815	ng	99
69) Atrazine	11.051	200	523276	63.950	ng	99
70) Pentachlorophenol	11.151	266	736673	106.719	ng	100
71) Phenanthrene	11.375	178	2438164	49.805	ng	99
72) Anthracene	11.427	178	2458121	51.244	ng	99
73) Carbazole	11.575	167	2220304	48.149	ng	100
74) Di-n-butylphthalate	11.910	149	2577929	49.151	ng	100
75) Fluoranthene	12.557	202	2511239	47.298	ng	100
77) Benzidine	12.680	184	795068	63.216	ng	100
78) Pyrene	12.786	202	2546390	48.925	ng	100
80) Butylbenzylphthalate	13.410	149	983900	51.771	ng	100
81) Benzo(a)anthracene	13.980	228	1908968	46.026	ng	100
82) 3,3'-Dichlorobenzidine	13.939	252	445199	35.727	ng	100
83) Chrysene	14.016	228	1790682	47.894	ng	100
84) Bis(2-ethylhexyl)phtha...	13.974	149	1158907	47.010	ng	99
85) Di-n-octyl phthalate	14.604	149	1808132	50.676	ng	97
87) Indeno(1,2,3-cd)pyrene	16.915	276	1936057	55.758	ng	98
88) Benzo(b)fluoranthene	15.033	252	1875492	48.792	ng	99
89) Benzo(k)fluoranthene	15.063	252	1446451	48.013	ng	99
90) Benzo(a)pyrene	15.398	252	1548930	51.917	ng	98
91) Dibenzo(a,h)anthracene	16.939	278	1564146	55.572	ng	98
92) Benzo(g,h,i)perylene	17.357	276	1473013	50.458	ng	98

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

