

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061825\
 Data File : BF142775.D
 Acq On : 18 Jun 2025 13:16
 Operator : RC/JU
 Sample : PB168505BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB168505BS

Quant Time: Jun 18 14:29:21 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF061125.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 11 05:56:09 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc Units	Dev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	6.886	152	81144	20.000 ng	0.00
21) Naphthalene-d8	8.169	136	313547	20.000 ng	-0.01
39) Acenaphthene-d10	9.927	164	171493	20.000 ng	-0.01
64) Phenanthrene-d10	11.416	188	297764	20.000 ng	0.00
76) Chrysene-d12	14.063	240	180043	20.000 ng	0.00
86) Perylene-d12	15.551	264	177074	20.000 ng	-0.01
System Monitoring Compounds					
5) 2-Fluorophenol	5.516	112	642728	134.903 ng	0.01
7) Phenol-d6	6.516	99	771821	137.654 ng	0.00
23) Nitrobenzene-d5	7.451	82	480756	83.913 ng	0.00
42) 2,4,6-Tribromophenol	10.722	330	252842	134.523 ng	0.00
45) 2-Fluorobiphenyl	9.245	172	1056545	81.851 ng	-0.01
79) Terphenyl-d14	13.004	244	1040355	79.362 ng	0.00
Target Compounds					
2) 1,4-Dioxane	2.740	88	74572	39.550 ng	98
3) Pyridine	3.498	79	207094	43.804 ng	97
4) n-Nitrosodimethylamine	3.451	42	119981	49.601 ng	98
6) Aniline	6.545	93	268193	35.738 ng	# 88
8) 2-Chlorophenol	6.669	128	251296	48.265 ng	100
9) Benzaldehyde	6.434	77	160795	46.309 ng	99
10) Phenol	6.534	94	294467	47.248 ng	91
11) bis(2-Chloroethyl)ether	6.622	93	219469	47.146 ng	99
12) 1,3-Dichlorobenzene	6.828	146	269109	45.290 ng	99
13) 1,4-Dichlorobenzene	6.904	146	275083	45.809 ng	100
14) 1,2-Dichlorobenzene	7.051	146	262483	45.600 ng	99
15) Benzyl Alcohol	7.028	79	205058	48.387 ng	100
16) 2,2'-oxybis(1-Chloropr...	7.157	45	347100	47.358 ng	97
17) 2-Methylphenol	7.139	107	193901	48.579 ng	99
18) Hexachloroethane	7.398	117	96483	44.834 ng	97
19) n-Nitroso-di-n-propyla...	7.298	70	165914	46.481 ng	99
20) 3+4-Methylphenols	7.292	107	241587	48.009 ng	97
22) Acetophenone	7.292	105	324787	46.567 ng	95
24) Nitrobenzene	7.469	77	243375	47.863 ng	98
25) Isophorone	7.710	82	445423	45.918 ng	99
26) 2-Nitrophenol	7.786	139	135453	47.732 ng	97
27) 2,4-Dimethylphenol	7.822	122	231910	47.999 ng	99
28) bis(2-Chloroethoxy)met...	7.916	93	280246	46.532 ng	99
29) 2,4-Dichlorophenol	8.028	162	211297	47.388 ng	98
30) 1,2,4-Trichlorobenzene	8.110	180	228090	46.293 ng	100
31) Naphthalene	8.192	128	725594	46.792 ng	100
32) Benzoic acid	7.945	122	133411	49.030 ng	100
33) 4-Chloroaniline	8.239	127	141096	22.662 ng	98
34) Hexachlorobutadiene	8.304	225	140810	44.847 ng	100
35) Caprolactam	8.610	113	66251m	55.044 ng	
36) 4-Chloro-3-methylphenol	8.722	107	222768	48.048 ng	99
37) 2-Methylnaphthalene	8.880	142	453908	46.246 ng	99
38) 1-Methylnaphthalene	8.980	142	471310	46.353 ng	99
40) 1,2,4,5-Tetrachloroben...	9.051	216	228887	46.114 ng	99
41) Hexachlorocyclopentadiene	9.033	237	286982	90.085 ng	99
43) 2,4,6-Trichlorophenol	9.163	196	155222	48.311 ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.204	196	166627	48.016	ng	99
46) 1,1'-Biphenyl	9.345	154	626959	47.081	ng	99
47) 2-Chloronaphthalene	9.375	162	460694	46.958	ng	99
48) 2-Nitroaniline	9.469	65	139312	49.926	ng	99
49) Acenaphthylene	9.792	152	786712	47.558	ng	100
50) Dimethylphthalate	9.651	163	546748	47.743	ng	100
51) 2,6-Dinitrotoluene	9.710	165	120652	48.747	ng	92
52) Acenaphthene	9.963	154	536473	52.306	ng	100
53) 3-Nitroaniline	9.880	138	73550	27.397	ng	100
54) 2,4-Dinitrophenol	9.992	184	143372	105.748	ng	91
55) Dibenzofuran	10.133	168	688889	47.168	ng	99
56) 4-Nitrophenol	10.045	139	196655	108.007	ng	95
57) 2,4-Dinitrotoluene	10.116	165	166124	50.764	ng	95
58) Fluorene	10.480	166	542310	47.021	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.251	232	137884	47.228	ng	98
60) Diethylphthalate	10.351	149	548034	48.261	ng	100
61) 4-Chlorophenyl-phenylether	10.469	204	258555	45.904	ng	99
62) 4-Nitroaniline	10.498	138	126267	52.574	ng	100
63) Azobenzene	10.627	77	475219	47.917	ng	100
65) 4,6-Dinitro-2-methylphthalate	10.527	198	89835	48.178	ng	95
66) n-Nitrosodiphenylamine	10.586	169	476429	46.548	ng	99
67) 4-Bromophenyl-phenylether	10.957	248	162420	46.211	ng	97
68) Hexachlorobenzene	11.027	284	178602	45.778	ng	98
69) Atrazine	11.116	200	150200	55.042	ng	99
70) Pentachlorophenol	11.222	266	199885	97.737	ng	99
71) Phenanthrene	11.439	178	763726	47.875	ng	100
72) Anthracene	11.492	178	772918	46.837	ng	99
73) Carbazole	11.645	167	704704	51.010	ng	100
74) Di-n-butylphthalate	11.974	149	803587	52.092	ng	100
75) Fluoranthene	12.633	202	762235	50.250	ng	99
77) Benzidine	12.751	184	299880	46.769	ng	99
78) Pyrene	12.863	202	764552	45.696	ng	100
80) Butylbenzylphthalate	13.474	149	277467	56.080	ng	99
81) Benzo(a)anthracene	14.051	228	590561	49.499	ng	99
82) 3,3'-Dichlorobenzidine	14.010	252	81013	21.057	ng	98
83) Chrysene	14.086	228	533597	48.441	ng	100
84) Bis(2-ethylhexyl)phthalate	14.033	149	364277	49.059	ng	100
85) Di-n-octyl phthalate	14.651	149	602523	42.295	ng	99
87) Indeno(1,2,3-cd)pyrene	17.068	276	599869	45.714	ng	99
88) Benzo(b)fluoranthene	15.115	252	521786	50.045	ng	99
89) Benzo(k)fluoranthene	15.145	252	490802	48.516	ng	100
90) Benzo(a)pyrene	15.492	252	492846	49.715	ng	99
91) Dibenzo(a,h)anthracene	17.086	278	494234	46.127	ng	99
92) Benzo(g,h,i)perylene	17.527	276	486563	45.667	ng	98

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

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