

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF100225\
 Data File : BF143864.D
 Acq On : 02 Oct 2025 19:38
 Operator : RC/JU
 Sample : Q3266-01MS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ENV-3-6FTMS

Quant Time: Oct 03 01:27:21 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF092325.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 24 05:00:00 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.887	152	119401	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	452831	20.000	ng	0.00
39) Acenaphthene-d10	9.922	164	241644	20.000	ng	0.00
64) Phenanthrene-d10	11.410	188	394825	20.000	ng	0.00
76) Chrysene-d12	14.057	240	306082	20.000	ng	0.00
86) Perylene-d12	15.533	264	431989	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	604335	82.232	ng	0.00
7) Phenol-d6	6.504	99	732194	86.227	ng	0.00
23) Nitrobenzene-d5	7.445	82	445326	60.197	ng	0.00
42) 2,4,6-Tribromophenol	10.716	330	195734	54.706	ng	0.00
45) 2-Fluorobiphenyl	9.239	172	781050	50.439	ng	0.00
79) Terphenyl-d14	12.998	244	1046086	69.507	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.711	88	145336	41.317	ng	92
3) Pyridine	3.469	79	394892	41.589	ng	87
4) n-Nitrosodimethylamine	3.422	42	202826	42.601	ng	84
6) Aniline	6.546	93	382496	31.036	ng	99
8) 2-Chlorophenol	6.669	128	360046	47.856	ng	96
9) Benzaldehyde	6.434	77	211201	28.402	ng	97
10) Phenol	6.522	94	470637	47.683	ng	99
11) bis(2-Chloroethyl)ether	6.616	93	351252	45.400	ng	97
12) 1,3-Dichlorobenzene	6.828	146	407642	46.666	ng	98
13) 1,4-Dichlorobenzene	6.904	146	408635	46.875	ng	99
14) 1,2-Dichlorobenzene	7.057	146	392158	47.497	ng	100
15) Benzyl Alcohol	7.022	79	322083	52.494	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.157	45	718122	39.143	ng	95
17) 2-Methylphenol	7.134	107	284360	48.820	ng	99
18) Hexachloroethane	7.398	117	142081	46.344	ng	97
19) n-Nitroso-di-n-propyla...	7.293	70	253743	46.469	ng	90
20) 3+4-Methylphenols	7.287	107	333283	46.985	ng	# 73
22) Acetophenone	7.293	105	499175	50.189	ng	94
24) Nitrobenzene	7.463	77	392104	47.818	ng	97
25) Isophorone	7.704	82	710337	49.274	ng	99
26) 2-Nitrophenol	7.781	139	184449	54.406	ng	94
27) 2,4-Dimethylphenol	7.816	122	336616	52.242	ng	99
28) bis(2-Chloroethoxy)met...	7.916	93	430461	45.960	ng	99
29) 2,4-Dichlorophenol	8.022	162	308452	48.651	ng	100
30) 1,2,4-Trichlorobenzene	8.104	180	325230	49.430	ng	98
31) Naphthalene	8.192	128	986012	45.485	ng	100
32) Benzoic acid	7.928	122	234383	75.394	ng	95
33) 4-Chloroaniline	8.234	127	115624	15.178	ng	97
34) Hexachlorobutadiene	8.304	225	229733	46.838	ng	99
35) Caprolactam	8.604	113	98125	60.552	ng	# 77
36) 4-Chloro-3-methylphenol	8.716	107	306778	51.593	ng	99
37) 2-Methylnaphthalene	8.881	142	612475	44.309	ng	100
38) 1-Methylnaphthalene	8.981	142	620851	46.926	ng	98
40) 1,2,4,5-Tetrachloroben...	9.045	216	356950	48.948	ng	99
41) Hexachlorocyclopentadiene	9.034	237	485641	105.078	ng	98
43) 2,4,6-Trichlorophenol	9.157	196	227991	49.930	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.192	196	234672	49.134	ng	100
46) 1,1'-Biphenyl	9.345	154	865505	49.480	ng	100
47) 2-Chloronaphthalene	9.369	162	629309	44.695	ng	100
48) 2-Nitroaniline	9.463	65	238647	56.079	ng	97
49) Acenaphthylene	9.787	152	1023535	52.805	ng	100
50) Dimethylphthalate	9.645	163	731465	49.664	ng	99
51) 2,6-Dinitrotoluene	9.704	165	156423	64.642	ng	94
52) Acenaphthene	9.957	154	622293	48.786	ng	99
53) 3-Nitroaniline	9.875	138	91487	31.191	ng	98
54) 2,4-Dinitrophenol	9.981	184	121210	100.703	ng	# 18
55) Dibenzofuran	10.134	168	848921	47.064	ng	100
56) 4-Nitrophenol	10.028	139	236657	100.124	ng	95
57) 2,4-Dinitrotoluene	10.110	165	212438	58.468	ng	92
58) Fluorene	10.475	166	641919	47.871	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.245	232	227326	55.559	ng	100
60) Diethylphthalate	10.345	149	698970	50.698	ng	99
61) 4-Chlorophenyl-phenyle...	10.463	204	330127	48.317	ng	98
62) 4-Nitroaniline	10.486	138	158530	57.481	ng	98
63) Azobenzene	10.628	77	737403	50.750	ng	93
65) 4,6-Dinitro-2-methylph...	10.516	198	95114	61.024	ng	# 81
66) n-Nitrosodiphenylamine	10.581	169	613626	47.258	ng	99
67) 4-Bromophenyl-phenylether	10.957	248	310198	46.102	ng	100
68) Hexachlorobenzene	11.022	284	374735	40.413	ng	93
69) Atrazine	11.110	200	197683	52.574	ng	98
70) Pentachlorophenol	11.216	266	416446	90.225	ng	99
71) Phenanthrene	11.439	178	940959	46.421	ng	99
72) Anthracene	11.492	178	966261	47.523	ng	100
73) Carbazole	11.645	167	836065	47.030	ng	100
74) Di-n-butylphthalate	11.975	149	1014489	49.620	ng	100
75) Fluoranthene	12.628	202	951417	46.667	ng	99
77) Benzidine	12.745	184	241360	32.360	ng	99
78) Pyrene	12.857	202	943184	62.677	ng	100
80) Butylbenzylphthalate	13.475	149	334241	64.200	ng	96
81) Benzo(a)anthracene	14.045	228	823820	50.101	ng	99
82) 3,3'-Dichlorobenzidine	14.004	252	159637	22.410	ng	99
83) Chrysene	14.080	228	760164	51.090	ng	99
84) Bis(2-ethylhexyl)phtha...	14.033	149	393368	51.547	ng	98
85) Di-n-octyl phthalate	14.645	149	673377	48.645	ng	# 93
87) Indeno(1,2,3-cd)pyrene	17.033	276	1617559	49.351	ng	97
88) Benzo(b)fluoranthene	15.098	252	1152982	52.411	ng	99
89) Benzo(k)fluoranthene	15.127	252	967267	43.764	ng	100
90) Benzo(a)pyrene	15.468	252	1075170	51.994	ng	99
91) Dibenzo(a,h)anthracene	17.051	278	1280220	47.754	ng	99
92) Benzo(g,h,i)perylene	17.486	276	1319988	51.078	ng	98

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

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