

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071625\
 Data File : BF143108.D
 Acq On : 16 Jul 2025 12:07
 Operator : RC/JU
 Sample : Q2126-08
 Misc : LOQ-WATER-5 PPM
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOQ-WATER-02-QT2-2025

Quant Time: Jul 16 12:28:02 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF071525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 15 17:53:25 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.969	152	141335	20.000	ng	0.00
21) Naphthalene-d8	8.245	136	536916	20.000	ng	0.00
39) Acenaphthene-d10	10.004	164	272300	20.000	ng	0.00
64) Phenanthrene-d10	11.492	188	444663	20.000	ng	0.00
76) Chrysene-d12	14.127	240	235543	20.000	ng	0.00
86) Perylene-d12	15.633	264	227368	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.592	112	939996	105.079	ng	0.00
7) Phenol-d6	6.587	99	1206482	107.294	ng	0.00
23) Nitrobenzene-d5	7.522	82	755341	78.920	ng	0.00
42) 2,4,6-Tribromophenol	10.792	330	297839	122.670	ng	0.00
45) 2-Fluorobiphenyl	9.322	172	1388592	67.787	ng	0.00
79) Terphenyl-d14	13.074	244	1208673	75.012	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.816	88	18349	4.121	ng	100
3) Pyridine	3.622	79	44147	3.915	ng	99
6) Aniline	6.628	93	68173	4.305	ng	98
8) 2-Chlorophenol	6.751	128	40446	4.500	ng	# 76
10) Phenol	6.598	94	56956	4.679	ng	# 68
11) bis(2-Chloroethyl)ether	6.698	93	41636	4.418	ng	98
12) 1,3-Dichlorobenzene	6.910	146	45360	4.467	ng	99
13) 1,4-Dichlorobenzene	6.987	146	45568	4.461	ng	99
14) 1,2-Dichlorobenzene	7.139	146	42740	4.387	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.239	45	76054	4.378	ng	100
17) 2-Methylphenol	7.198	107	32558	4.198	ng	95
18) Hexachloroethane	7.486	117	14626	4.398	ng	99
19) n-Nitroso-di-n-propyla...	7.363	70	30839	4.477	ng	95
22) Acetophenone	7.363	105	60332	4.661	ng	95
24) Nitrobenzene	7.539	77	45597	4.935	ng	99
25) Isophorone	7.775	82	79829	4.315	ng	99
26) 2-Nitrophenol	7.857	139	11850	6.877	ng	95
27) 2,4-Dimethylphenol	7.886	122	38952	4.537	ng	98
28) bis(2-Chloroethoxy)met...	7.986	93	50369	4.490	ng	97
29) 2,4-Dichlorophenol	8.092	162	30459	4.336	ng	98
30) 1,2,4-Trichlorobenzene	8.186	180	35933	4.507	ng	99
31) Naphthalene	8.269	128	123701	4.640	ng	99
33) 4-Chloroaniline	8.310	127	46601	4.346	ng	98
34) Hexachlorobutadiene	8.392	225	21218	4.450	ng	98
36) 4-Chloro-3-methylphenol	8.775	107	33020	4.220	ng	98
37) 2-Methylnaphthalene	8.957	142	72600	4.564	ng	98
38) 1-Methylnaphthalene	9.057	142	77200	4.680	ng	99
40) 1,2,4,5-Tetrachloroben...	9.128	216	36664	4.620	ng	98
43) 2,4,6-Trichlorophenol	9.228	196	19829	4.022	ng	97
44) 2,4,5-Trichlorophenol	9.263	196	21550	4.168	ng	97
46) 1,1'-Biphenyl	9.422	154	99699	4.591	ng	99
47) 2-Chloronaphthalene	9.445	162	71802	4.501	ng	99
48) 2-Nitroaniline	9.528	65	14751	6.068	ng	95
49) Acenaphthylene	9.863	152	118162	4.445	ng	99
50) Dimethylphthalate	9.710	163	77182	4.505	ng	99
51) 2,6-Dinitrotoluene	9.769	165	11837	6.089	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
52) Acenaphthene	10.033	154	70980	4.525	ng	99
53) 3-Nitroaniline	9.933	138	14209	3.839	ng	# 93
55) Dibenzofuran	10.204	168	108220	4.628	ng	97
57) 2,4-Dinitrotoluene	10.169	165	12579	6.164	ng	96
58) Fluorene	10.551	166	82319	4.696	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.322	232	17445	4.246	ng	98
60) Diethylphthalate	10.416	149	71822	4.346	ng	100
61) 4-Chlorophenyl-phenyle...	10.539	204	38270	4.555	ng	96
62) 4-Nitroaniline	10.545	138	12756	3.963	ng	96
63) Azobenzene	10.704	77	79730	4.362	ng	99
66) n-Nitrosodiphenylamine	10.657	169	68367	4.377	ng	99
67) 4-Bromophenyl-phenylether	11.033	248	21389	4.230	ng	96
68) Hexachlorobenzene	11.104	284	23075	4.301	ng	98
69) Atrazine	11.174	200	15480	3.874	ng	100
71) Phenanthrene	11.510	178	109851	4.578	ng	100
72) Anthracene	11.563	178	106216	4.401	ng	99
73) Carbazole	11.710	167	95358	4.411	ng	99
74) Di-n-butylphthalate	12.051	149	73329	3.695	ng	99
75) Fluoranthene	12.704	202	92470	4.238	ng	99
78) Pyrene	12.933	202	97548	4.464	ng	99
80) Butylbenzylphthalate	13.545	149	11270	5.796	ng	92
81) Benzo(a)anthracene	14.115	228	63108	4.021	ng	98
83) Chrysene	14.151	228	63069	4.363	ng	99
84) Bis(2-ethylhexyl)phtha...	14.110	149	17064	5.805	ng	99
87) Indeno(1,2,3-cd)pyrene	17.168	276	68340	4.041	ng	99
88) Benzo(b)fluoranthene	15.186	252	51884	3.956	ng	100
89) Benzo(k)fluoranthene	15.215	252	47857	3.825	ng	98
90) Benzo(a)pyrene	15.568	252	45828	3.695	ng	98
91) Dibenzo(a,h)anthracene	17.186	278	56845	4.112	ng	97
92) Benzo(g,h,i)perylene	17.627	276	57030	4.046	ng	96

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

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