

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112025\
 Data File : BF144303.D
 Acq On : 20 Nov 2025 19:26
 Operator : RC/JU
 Sample : Q3662-02MS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 B3-0.0-1.0-20251117MS

Quant Time: Nov 20 23:59:36 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 05 18:37:33 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.863	152	47796	20.000	ng	-0.01
21) Naphthalene-d8	8.145	136	175247	20.000	ng	-0.01
39) Acenaphthene-d10	9.904	164	88326	20.000	ng	-0.01
64) Phenanthrene-d10	11.392	188	153411	20.000	ng	-0.01
76) Chrysene-d12	14.045	240	129309	20.000	ng	0.00
86) Perylene-d12	15.521	264	103273	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.493	112	215853	70.672	ng	0.00
7) Phenol-d6	6.498	99	242523	70.078	ng	0.00
23) Nitrobenzene-d5	7.428	82	148501	48.940	ng	0.00
42) 2,4,6-Tribromophenol	10.698	330	72543	65.449	ng	0.00
45) 2-Fluorobiphenyl	9.222	172	301591	50.430	ng	0.00
79) Terphenyl-d14	12.980	244	342368	42.056	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	2.669	88	63317	50.143	ng	98
3) Pyridine	3.434	79	164219	46.615	ng	97
4) n-Nitrosodimethylamine	3.381	42	90534	49.199	ng	91
6) Aniline	6.528	93	67950	13.781	ng	# 85
8) 2-Chlorophenol	6.651	128	135084	45.098	ng	98
9) Benzaldehyde	6.416	77	73252	37.484	ng	98
10) Phenol	6.510	94	184062	43.531	ng	96
11) bis(2-Chloroethyl)ether	6.598	93	144683	46.959	ng	98
12) 1,3-Dichlorobenzene	6.804	146	171695	46.453	ng	98
13) 1,4-Dichlorobenzene	6.881	146	172341	46.542	ng	98
14) 1,2-Dichlorobenzene	7.034	146	165526	46.637	ng	98
15) Benzyl Alcohol	7.004	79	126922	46.316	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.139	45	260222	45.864	ng	99
17) 2-Methylphenol	7.122	107	107301	45.030	ng	98
18) Hexachloroethane	7.375	117	52802	42.920	ng	97
19) n-Nitroso-di-n-propyla...	7.275	70	102677	45.068	ng	98
20) 3+4-Methylphenols	7.275	107	125726	44.758	ng	# 89
22) Acetophenone	7.275	105	178429	47.462	ng	97
24) Nitrobenzene	7.445	77	151066	45.853	ng	99
25) Isophorone	7.686	82	267312	46.574	ng	99
26) 2-Nitrophenol	7.763	139	63205	38.755	ng	97
27) 2,4-Dimethylphenol	7.798	122	122563	50.140	ng	99
28) bis(2-Chloroethoxy)met...	7.892	93	167606	46.764	ng	97
29) 2,4-Dichlorophenol	8.010	162	120659	42.827	ng	98
30) 1,2,4-Trichlorobenzene	8.086	180	133419	46.118	ng	96
31) Naphthalene	8.169	128	374098	44.876	ng	99
32) Benzoic acid	7.910	122	60596	33.819	ng	96
33) 4-Chloroaniline	8.222	127	34485	11.342	ng	98
34) Hexachlorobutadiene	8.281	225	91710	42.063	ng	97
35) Caprolactam	8.586	113	34548	44.759	ng	93
36) 4-Chloro-3-methylphenol	8.704	107	104964	41.571	ng	99
37) 2-Methylnaphthalene	8.857	142	224064	40.201	ng	98
38) 1-Methylnaphthalene	8.957	142	220613	41.022	ng	100
40) 1,2,4,5-Tetrachloroben...	9.028	216	134695	46.531	ng	99
41) Hexachlorocyclopentadiene	9.010	237	24125	29.996	ng	96
43) 2,4,6-Trichlorophenol	9.139	196	87192	44.255	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.186	196	92551	43.586	ng	99
46) 1,1'-Biphenyl	9.322	154	290901	47.183	ng	100
47) 2-Chloronaphthalene	9.351	162	237319	45.125	ng	99
48) 2-Nitroaniline	9.445	65	75092	49.486	ng	99
49) Acenaphthylene	9.763	152	364796	50.901	ng	100
50) Dimethylphthalate	9.622	163	282211	48.229	ng	100
51) 2,6-Dinitrotoluene	9.686	165	57819	48.812	ng	95
52) Acenaphthene	9.939	154	194937	40.675	ng	99
53) 3-Nitroaniline	9.857	138	43949	33.941	ng	100
54) 2,4-Dinitrophenol	9.975	184	5029	7.031	ng	93
55) Dibenzofuran	10.110	168	301970	44.988	ng	100
56) 4-Nitrophenol	10.028	139	84871	90.611	ng	94
57) 2,4-Dinitrotoluene	10.092	165	73241	46.188	ng	# 99
58) Fluorene	10.451	166	234298	45.911	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.233	232	70902	47.194	ng	100
60) Diethylphthalate	10.322	149	256137	47.485	ng	99
61) 4-Chlorophenyl-phenylether	10.445	204	126225	43.421	ng	99
62) 4-Nitroaniline	10.475	138	58579	47.505	ng	95
63) Azobenzene	10.604	77	246183	49.039	ng	95
65) 4,6-Dinitro-2-methylph...	10.504	198	5944	5.703	ng	98
66) n-Nitrosodiphenylamine	10.563	169	225030	45.607	ng	100
67) 4-Bromophenyl-phenylether	10.933	248	82595	42.180	ng	97
68) Hexachlorobenzene	11.004	284	99968	43.343	ng	96
69) Atrazine	11.092	200	75460	48.090	ng	98
70) Pentachlorophenol	11.204	266	97266	84.688	ng	97
71) Phenanthrene	11.422	178	364570	47.730	ng	99
72) Anthracene	11.469	178	364256	46.900	ng	100
73) Carbazole	11.627	167	333099	50.437	ng	99
74) Di-n-butylphthalate	11.951	149	392037	49.104	ng	99
75) Fluoranthene	12.610	202	455093	57.678	ng	99
77) Benzidine	12.733	184	133834	34.990	ng	97
78) Pyrene	12.845	202	471853	45.257	ng	100
80) Butylbenzylphthalate	13.457	149	177107	49.012	ng	97
81) Benzo(a)anthracene	14.033	228	436124	48.663	ng	99
82) 3,3'-Dichlorobenzidine	13.992	252	105242	34.259	ng	97
83) Chrysene	14.068	228	402572	48.660	ng	99
84) Bis(2-ethylhexyl)phtha...	14.015	149	247104	49.953	ng	97
85) Di-n-octyl phthalate	14.627	149	409575	47.989	ng	100
87) Indeno(1,2,3-cd)pyrene	17.021	276	280692	37.863	ng	98
88) Benzo(b)fluoranthene	15.092	252	338864	57.716	ng	99
89) Benzo(k)fluoranthene	15.115	252	292518	53.229	ng	99
90) Benzo(a)pyrene	15.462	252	276440	51.037	ng	97
91) Dibenzo(a,h)anthracene	17.033	278	223389	37.521	ng	98
92) Benzo(g,h,i)perylene	17.474	276	218137	37.102	ng	99

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

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