

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF091025\
 Data File : BF143688.D
 Acq On : 09 Sep 2025 16:59
 Operator : RC/JU
 Sample : Q2893-01
 Misc : LOD-MDL-SOIL 8 PPM
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOD-MDL-SOIL-01-QT3-2025

Quant Time: Sep 09 17:50:04 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF091025.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 09 15:20:16 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.887	152	45543	20.000	ng	0.00
21) Naphthalene-d8	8.169	136	173489	20.000	ng	0.00
39) Acenaphthene-d10	9.922	164	102171	20.000	ng	0.00
64) Phenanthrene-d10	11.410	188	173016	20.000	ng	0.00
76) Chrysene-d12	14.045	240	103451	20.000	ng	0.00
86) Perylene-d12	15.516	264	105156	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	353441	132.407	ng	0.00
7) Phenol-d6	6.510	99	442433	136.124	ng	0.00
23) Nitrobenzene-d5	7.451	82	277825	99.608	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	143175	139.052	ng	0.00
45) 2-Fluorobiphenyl	9.245	172	623246	92.148	ng	0.00
79) Terphenyl-d14	12.998	244	657958	103.067	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.681	88	9964	8.368	ng	96
3) Pyridine	3.475	79	24022	7.810	ng	96
4) n-Nitrosodimethylamine	3.375	42	9241	7.909	ng	96
6) Aniline	6.551	93	36335	8.441	ng	97
8) 2-Chlorophenol	6.675	128	24031	8.335	ng	97
9) Benzaldehyde	6.440	77	18761	7.154	ng	96
10) Phenol	6.522	94	28840	8.027	ng	81
11) bis(2-Chloroethyl)ether	6.622	93	22096	8.235	ng	98
12) 1,3-Dichlorobenzene	6.828	146	28965	8.785	ng	98
13) 1,4-Dichlorobenzene	6.904	146	28648	8.608	ng	99
14) 1,2-Dichlorobenzene	7.057	146	26496	8.431	ng	97
15) Benzyl Alcohol	7.022	79	19612	8.295	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.163	45	28383	8.085	ng	97
17) 2-Methylphenol	7.122	107	19146	8.256	ng	99
18) Hexachloroethane	7.404	117	9066	8.433	ng	99
19) n-Nitroso-di-n-propyla...	7.293	70	16995	8.514	ng	99
20) 3+4-Methylphenols	7.281	107	24939	8.166	ng	# 87
22) Acetophenone	7.293	105	34917	8.911	ng	98
24) Nitrobenzene	7.463	77	25091	8.676	ng	97
25) Isophorone	7.698	82	45488	8.505	ng	100
26) 2-Nitrophenol	7.781	139	11325	7.964	ng	97
27) 2,4-Dimethylphenol	7.810	122	23307	9.130	ng	99
28) bis(2-Chloroethoxy)met...	7.910	93	28316	8.615	ng	99
29) 2,4-Dichlorophenol	8.016	162	21194	8.254	ng	99
30) 1,2,4-Trichlorobenzene	8.110	180	23225	8.730	ng	99
31) Naphthalene	8.193	128	72225	8.410	ng	99
32) Benzoic acid	7.857	122	8694	5.582	ng	97
33) 4-Chloroaniline	8.234	127	28482	9.526	ng	96
34) Hexachlorobutadiene	8.310	225	14438	8.321	ng	97
35) Caprolactam	8.563	113	5535	8.099	ng	96
36) 4-Chloro-3-methylphenol	8.698	107	21154	8.355	ng	98
37) 2-Methylnaphthalene	8.881	142	45032	7.660	ng	99
38) 1-Methylnaphthalene	8.981	142	46613	8.262	ng	97
40) 1,2,4,5-Tetrachloroben...	9.045	216	24539	8.389	ng	98
41) Hexachlorocyclopentadiene	9.034	237	14591	9.716	ng	97
43) 2,4,6-Trichlorophenol	9.151	196	15537	7.955	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.187	196	16443	8.281	ng	99
46) 1,1'-Biphenyl	9.345	154	64249	8.735	ng	100
47) 2-Chloronaphthalene	9.369	162	47633	8.260	ng	98
48) 2-Nitroaniline	9.457	65	12281	8.048	ng	96
49) Acenaphthylene	9.781	152	76954	9.316	ng	100
50) Dimethylphthalate	9.640	163	56349	8.477	ng	99
51) 2,6-Dinitrotoluene	9.698	165	10999	8.654	ng	97
52) Acenaphthene	9.957	154	46463	8.163	ng	100
53) 3-Nitroaniline	9.863	138	11371	8.313	ng	89
54) 2,4-Dinitrophenol	9.969	184	3429	10.259	ng	# 1
55) Dibenzofuran	10.128	168	69751	8.522	ng	98
56) 4-Nitrophenol	10.004	139	8286	7.615	ng	96
57) 2,4-Dinitrotoluene	10.098	165	14024	8.107	ng	# 98
58) Fluorene	10.469	166	55329	8.636	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.239	232	14186	8.811	ng	# 97
60) Diethylphthalate	10.339	149	52683	8.442	ng	99
61) 4-Chlorophenyl-phenyle...	10.463	204	27490	8.460	ng	100
62) 4-Nitroaniline	10.469	138	10817	8.350	ng	97
63) Azobenzene	10.622	77	47209	8.556	ng	97
65) 4,6-Dinitro-2-methylph...	10.504	198	5511	6.256	ng	98
66) n-Nitrosodiphenylamine	10.581	169	47610	8.400	ng	99
67) 4-Bromophenyl-phenylether	10.951	248	17298	8.291	ng	95
68) Hexachlorobenzene	11.016	284	17593	7.986	ng	96
69) Atrazine	11.098	200	14106	7.937	ng	95
70) Pentachlorophenol	11.204	266	8766	6.582	ng	98
71) Phenanthrene	11.434	178	80496	8.444	ng	99
72) Anthracene	11.481	178	80940	8.444	ng	99
73) Carbazole	11.633	167	66847	8.515	ng	99
74) Di-n-butylphthalate	11.969	149	77926	8.245	ng	99
75) Fluoranthene	12.616	202	73412	8.389	ng	99
77) Benzidine	12.739	184	27459	8.140	ng	# 96
78) Pyrene	12.845	202	73617	8.900	ng	99
80) Butylbenzylphthalate	13.469	149	20832	8.038	ng	98
81) Benzo(a)anthracene	14.033	228	56450	8.482	ng	99
82) 3,3'-Dichlorobenzidine	13.992	252	17266	8.433	ng	99
83) Chrysene	14.069	228	49588	8.135	ng	96
84) Bis(2-ethylhexyl)phtha...	14.027	149	29801	8.059	ng	# 98
85) Di-n-octyl phthalate	14.639	149	46100	6.957	ng	99
87) Indeno(1,2,3-cd)pyrene	16.992	276	50917	7.095	ng	99
88) Benzo(b)fluoranthene	15.080	252	53022	8.470	ng	99
89) Benzo(k)fluoranthene	15.110	252	43120	7.633	ng	99
90) Benzo(a)pyrene	15.451	252	45182	8.216	ng	99
91) Dibenzo(a,h)anthracene	17.015	278	42471	7.133	ng	98
92) Benzo(g,h,i)perylene	17.439	276	40168	7.212	ng	97

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

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[illegible]