

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF121525\
 Data File : BF144470.D
 Acq On : 15 Dec 2025 13:12
 Operator : RC/JU
 Sample : Q3530-01
 Misc : LOD-SOIL-4NG
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 LOD-MDL-SOIL-01-QT4-2025

Quant Time: Dec 15 13:32:07 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF121225.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 12 15:57:03 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.793	152	264563	20.000	ng	0.00
21) Naphthalene-d8	8.075	136	1067346	20.000	ng	0.00
39) Acenaphthene-d10	9.828	164	578619	20.000	ng	0.00
64) Phenanthrene-d10	11.310	188	985264	20.000	ng	0.00
76) Chrysene-d12	13.939	240	798537	20.000	ng	0.00
86) Perylene-d12	15.374	264	615507	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.399	112	1803149	102.766	ng	0.00
7) Phenol-d6	6.416	99	2332562	106.579	ng	0.00
23) Nitrobenzene-d5	7.357	82	1557140	91.294	ng	0.00
42) 2,4,6-Tribromophenol	10.616	330	657646	125.714	ng	0.00
45) 2-Fluorobiphenyl	9.151	172	2677437	70.986	ng	0.00
79) Terphenyl-d14	12.898	244	3169271	62.660	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.493	88	28466	3.374	ng	97
3) Pyridine	3.287	79	74460	3.209	ng	98
4) n-Nitrosodimethylamine	3.158	42	35526	3.524	ng	# 96
6) Aniline	6.457	93	119006	3.898	ng	98
8) 2-Chlorophenol	6.575	128	74026	3.935	ng	# 76
9) Benzaldehyde	6.346	77	61478	3.588	ng	96
10) Phenol	6.434	94	96505	3.736	ng	79
11) bis(2-Chloroethyl)ether	6.528	93	73476	3.892	ng	99
12) 1,3-Dichlorobenzene	6.734	146	82342	4.051	ng	99
13) 1,4-Dichlorobenzene	6.810	146	82098	4.026	ng	99
14) 1,2-Dichlorobenzene	6.963	146	77210	3.987	ng	98
15) Benzyl Alcohol	6.928	79	61132	3.790	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.069	45	114474	3.593	ng	99
17) 2-Methylphenol	7.034	107	59406	3.746	ng	98
18) Hexachloroethane	7.310	117	26511	3.741	ng	97
19) n-Nitroso-di-n-propyla...	7.198	70	53281	3.854	ng	99
20) 3+4-Methylphenols	7.187	107	78394	4.024	ng	# 88
22) Acetophenone	7.198	105	108194	4.087	ng	95
24) Nitrobenzene	7.375	77	83105	4.491	ng	99
25) Isophorone	7.610	82	140315	3.645	ng	98
26) 2-Nitrophenol	7.687	139	25431	6.896	ng	98
27) 2,4-Dimethylphenol	7.722	122	65680	3.576	ng	95
28) bis(2-Chloroethoxy)met...	7.822	93	89463	3.772	ng	99
29) 2,4-Dichlorophenol	7.922	162	58507	3.834	ng	98
30) 1,2,4-Trichlorobenzene	8.016	180	68035	4.110	ng	98
31) Naphthalene	8.092	128	227015	3.952	ng	99
32) Benzoic acid	7.757	122	18349	7.981	ng	96
33) 4-Chloroaniline	8.140	127	91390	3.958	ng	98
34) Hexachlorobutadiene	8.216	225	39328	3.843	ng	98
35) Caprolactam	8.463	113	17022	3.378	ng	97
36) 4-Chloro-3-methylphenol	8.610	107	60871	3.674	ng	98
37) 2-Methylnaphthalene	8.787	142	133851	3.541	ng	99
38) 1-Methylnaphthalene	8.887	142	142721	3.878	ng	98
40) 1,2,4,5-Tetrachloroben...	8.951	216	67932	4.125	ng	97
41) Hexachlorocyclopentadiene	8.939	237	26684	2.738	ng	100
43) 2,4,6-Trichlorophenol	9.057	196	39097	3.739	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.092	196	41822	3.810	ng	97
46) 1,1'-Biphenyl	9.251	154	189961	4.201	ng	97
47) 2-Chloronaphthalene	9.275	162	138129	4.001	ng	97
48) 2-Nitroaniline	9.363	65	30267	6.458	ng	98
49) Acenaphthylene	9.687	152	215901	3.948	ng	98
50) Dimethylphthalate	9.545	163	149140	3.816	ng	99
51) 2,6-Dinitrotoluene	9.604	165	26345	7.287	ng	97
52) Acenaphthene	9.857	154	135704	3.980	ng	100
53) 3-Nitroaniline	9.769	138	31361	6.361	ng	# 95
54) 2,4-Dinitrophenol	9.875	184	5689	8.671	ng	# 36
55) Dibenzofuran	10.028	168	195363	4.026	ng	97
56) 4-Nitrophenol	9.916	139	21613	3.247	ng	96
57) 2,4-Dinitrotoluene	10.004	165	30795	6.781	ng	96
58) Fluorene	10.369	166	153829	4.187	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.145	232	35389	4.248	ng	97
60) Diethylphthalate	10.245	149	139587	3.768	ng	98
61) 4-Chlorophenyl-phenyle...	10.369	204	73317	4.067	ng	100
62) 4-Nitroaniline	10.375	138	30287	5.835	ng	99
63) Azobenzene	10.528	77	145397	3.714	ng	98
65) 4,6-Dinitro-2-methylph...	10.404	198	9088	8.672	ng	82
66) n-Nitrosodiphenylamine	10.481	169	128422	3.803	ng	99
67) 4-Bromophenyl-phenylether	10.857	248	42266	3.675	ng	98
68) Hexachlorobenzene	10.916	284	48618	3.829	ng	97
69) Atrazine	11.004	200	39003	3.688	ng	99
70) Pentachlorophenol	11.104	266	23298	3.217	ng	97
71) Phenanthrene	11.333	178	227327	4.002	ng	98
72) Anthracene	11.381	178	224411	3.904	ng	98
73) Carbazole	11.533	167	202475	4.057	ng	98
74) Di-n-butylphthalate	11.875	149	193709	3.631	ng	99
75) Fluoranthene	12.516	202	227616	4.058	ng	99
77) Benzidine	12.639	184	89353	3.143	ng	98
78) Pyrene	12.745	202	239560	3.274	ng	99
80) Butylbenzylphthalate	13.369	149	63487	5.702	ng	100
81) Benzo(a)anthracene	13.927	228	199434	3.546	ng	98
82) 3,3'-Dichlorobenzidine	13.892	252	51049	3.204	ng	99
83) Chrysene	13.963	228	196419	3.817	ng	99
84) Bis(2-ethylhexyl)phtha...	13.933	149	75581	5.427	ng	99
85) Di-n-octyl phthalate	14.545	149	84522	8.187	ng	99
87) Indeno(1,2,3-cd)pyrene	16.780	276	126383	2.861	ng	99
88) Benzo(b)fluoranthene	14.957	252	144871	3.585	ng	100
89) Benzo(k)fluoranthene	14.986	252	151961	3.977	ng	99
90) Benzo(a)pyrene	15.310	252	124565	3.527	ng	100
91) Dibenzo(a,h)anthracene	16.798	278	101756	2.852	ng	98
92) Benzo(g,h,i)perylene	17.204	276	104623	2.884	ng	99

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

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