

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM022825\
 Data File : BM049683.D
 Acq On : 28 Feb 2025 17:55
 Operator : RC/JU
 Sample : Q1458-01MS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SP-SOILMS

Quant Time: Feb 28 21:59:40 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM020525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 06 04:55:28 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.804	152	347305	20.000	ng	-0.02
21) Naphthalene-d8	10.598	136	1210606	20.000	ng	#-0.03
39) Acenaphthene-d10	14.445	164	762853	20.000	ng	-0.02
64) Phenanthrene-d10	17.186	188	1502819	20.000	ng	-0.02
76) Chrysene-d12	21.427	240	1693591	20.000	ng	-0.02
86) Perylene-d12	24.438	264	1738682	20.000	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	2490582	115.319	ng	0.00
7) Phenol-d6	6.981	99	3098432	109.504	ng	0.00
23) Nitrobenzene-d5	8.957	82	1787386	75.944	ng	-0.02
42) 2,4,6-Tribromophenol	15.933	330	1378582	152.512	ng	-0.02
45) 2-Fluorobiphenyl	13.069	172	4386643	74.493	ng	-0.02
79) Terphenyl-d14	19.809	244	6599150	59.502	ng	-0.02
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.281	88	372494	41.020	ng	91
3) Pyridine	3.681	79	909357	36.036	ng	90
4) n-Nitrosodimethylamine	3.587	42	367200	31.102	ng	# 77
6) Aniline	7.128	93	477327	17.799	ng	93
8) 2-Chlorophenol	7.375	128	927494	43.008	ng	93
9) Benzaldehyde	6.940	77	339702	22.674	ng	93
10) Phenol	7.010	94	1151957	41.334	ng	95
11) bis(2-Chloroethyl)ether	7.228	93	893658	39.636	ng	95
12) 1,3-Dichlorobenzene	7.692	146	1090697	43.572	ng	# 95
13) 1,4-Dichlorobenzene	7.839	146	1100624	43.469	ng	98
14) 1,2-Dichlorobenzene	8.157	146	1058311	43.289	ng	97
15) Benzyl Alcohol	8.045	79	746511	37.922	ng	95
16) 2,2'-oxybis(1-Chloropr...	8.328	45	1026803	33.830	ng	93
17) 2-Methylphenol	8.257	107	717422	42.100	ng	96
18) Hexachloroethane	8.886	117	383740	41.492	ng	91
19) n-Nitroso-di-n-propyla...	8.610	70	644418	34.779	ng	91
20) 3+4-Methylphenols	8.581	107	967373	43.003	ng	97
22) Acetophenone	8.622	105	1433077	43.951	ng	96
24) Nitrobenzene	9.004	77	904790	39.569	ng	92
25) Isophorone	9.534	82	1695898	39.955	ng	96
26) 2-Nitrophenol	9.710	139	458467	56.059	ng	90
27) 2,4-Dimethylphenol	9.781	122	739537	55.103	ng	96
28) bis(2-Chloroethoxy)met...	10.010	93	1094156	40.024	ng	97
29) 2,4-Dichlorophenol	10.251	162	924470	50.018	ng	95
30) 1,2,4-Trichlorobenzene	10.463	180	1036567	45.776	ng	99
31) Naphthalene	10.651	128	2736140	42.854	ng	99
32) Benzoic acid	9.916	122	540207	59.118	ng	93
33) 4-Chloroaniline	10.757	127	312299	13.068	ng	95
34) Hexachlorobutadiene	10.945	225	656379	48.404	ng	99
35) Caprolactam	11.539	113	270402	47.956	ng	# 80
36) 4-Chloro-3-methylphenol	11.898	107	786181	42.653	ng	92
37) 2-Methylnaphthalene	12.263	142	1882028	41.866	ng	99
38) 1-Methylnaphthalene	12.486	142	1836236	41.983	ng	99
40) 1,2,4,5-Tetrachloroben...	12.633	216	1349027	53.786	ng	99
41) Hexachlorocyclopentadiene	12.622	237	1201438	187.639	ng	99
43) 2,4,6-Trichlorophenol	12.880	196	754502	55.127	ng	98

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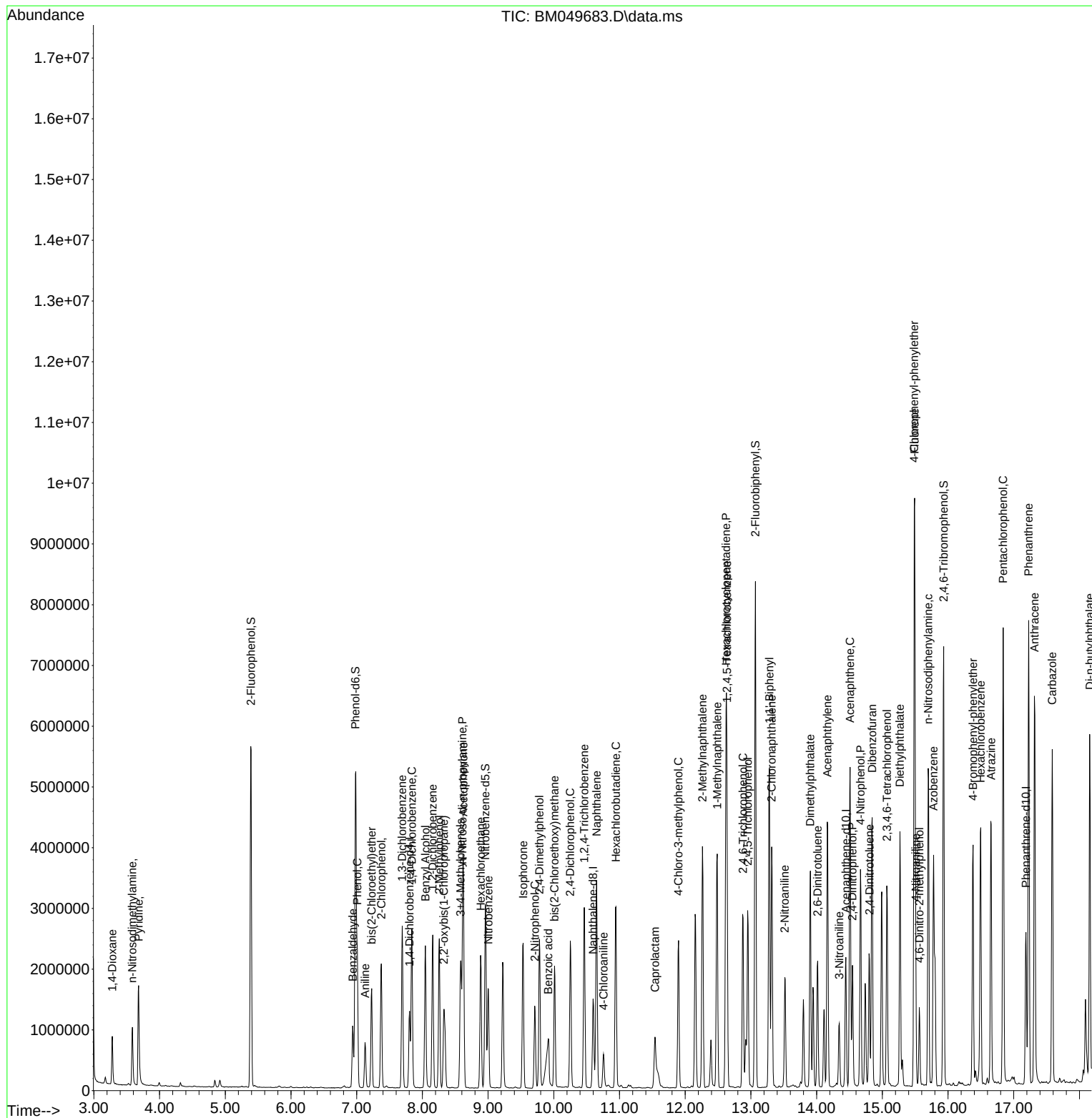
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.951	196	812939	54.049	ng	97
46) 1,1'-Biphenyl	13.280	154	2844593	48.881	ng	99
47) 2-Chloronaphthalene	13.316	162	2001974	44.463	ng	98
48) 2-Nitroaniline	13.516	65	463622	39.823	ng	88
49) Acenaphthylene	14.163	152	3119412	46.984	ng	99
50) Dimethylphthalate	13.904	163	2357564	43.034	ng	100
51) 2,6-Dinitrotoluene	14.016	165	488431	51.524	ng	# 82
52) Acenaphthene	14.510	154	2141511m	48.595	ng	
53) 3-Nitroaniline	14.345	138	273083	28.294	ng	# 90
54) 2,4-Dinitrophenol	14.545	184	454515	106.920	ng	# 66
55) Dibenzofuran	14.845	168	3010988	42.910	ng	96
56) 4-Nitrophenol	14.668	139	960944	118.109	ng	# 82
57) 2,4-Dinitrotoluene	14.804	165	688345	51.085	ng	# 78
58) Fluorene	15.492	166	2525629	42.074	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.068	232	684061	53.832	ng	92
60) Diethylphthalate	15.268	149	2273425	41.292	ng	98
61) 4-Chlorophenyl-phenyle...	15.486	204	1346603	42.147	ng	99
62) 4-Nitroaniline	15.510	138	470337	38.565	ng	92
63) Azobenzene	15.780	77	2490363	43.335	ng	90
65) 4,6-Dinitro-2-methylph...	15.563	198	303979	61.318	ng	85
66) n-Nitrosodiphenylamine	15.698	169	2093131	44.164	ng	99
67) 4-Bromophenyl-phenylether	16.380	248	785385	44.985	ng	95
68) Hexachlorobenzene	16.498	284	890427	44.436	ng	93
69) Atrazine	16.657	200	842821	59.823	ng	97
70) Pentachlorophenol	16.839	266	1311573	95.339	ng	99
71) Phenanthrene	17.227	178	4455337	52.039	ng	99
72) Anthracene	17.315	178	4080532	48.135	ng	100
73) Carbazole	17.586	167	3518716	47.292	ng	100
74) Di-n-butylphthalate	18.156	149	4099601	43.423	ng	99
75) Fluoranthene	19.245	202	6295201	62.805	ng	98
77) Benzidine	19.427	184	2046039	136.423	ng	100
78) Pyrene	19.603	202	6567296	51.788	ng	99
80) Butylbenzylphthalate	20.509	149	1870347	47.399	ng	# 85
81) Benzo(a)anthracene	21.409	228	5878271	51.214	ng	99
82) 3,3'-Dichlorobenzidine	21.327	252	1499205	48.540	ng	99
83) Chrysene	21.468	228	5335147	49.613	ng	99
84) Bis(2-ethylhexyl)phtha...	21.339	149	2892487	42.760	ng	# 97
85) Di-n-octyl phthalate	22.480	149	4923044	46.131	ng	# 92
87) Indeno(1,2,3-cd)pyrene	27.850	276	6376944	67.159	ng	# 96
88) Benzo(b)fluoranthene	23.491	252	5758593	48.145	ng	99
89) Benzo(k)fluoranthene	23.550	252	5331064	45.889	ng	99
90) Benzo(a)pyrene	24.303	252	5395741	56.924	ng	# 98
91) Dibenzo(a,h)anthracene	27.903	278	4922455	65.593	ng	98
92) Benzo(g,h,i)perylene	28.909	276	4941599	58.947	ng	97

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

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