

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082724\
 Data File : BM047361.D
 Acq On : 27 Aug 2024 13:17
 Operator : RC/JU
 Sample : P3726-01MSD 2X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 MOO-SOIL-PILEMSD

Quant Time: Aug 27 14:15:12 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM082624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 26 18:11:40 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.346	152	330759	20.000	ng	0.00
21) Naphthalene-d8	10.099	136	1270835	20.000	ng	0.00
39) Acenaphthene-d10	14.010	164	872733	20.000	ng	0.00
64) Phenanthrene-d10	16.775	188	1776554	20.000	ng	0.00
76) Chrysene-d12	21.027	240	1136290	20.000	ng	0.00
86) Perylene-d12	23.721	264	1306738	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.011	112	750409	43.527	ng	0.00
7) Phenol-d6	6.569	99	983254	41.867	ng	0.00
23) Nitrobenzene-d5	8.493	82	650417	28.398	ng	0.00
42) 2,4,6-Tribromophenol	15.522	330	450845	42.818	ng	0.00
45) 2-Fluorobiphenyl	12.610	172	1499578	26.401	ng	0.00
79) Terphenyl-d14	19.445	244	1839344	28.813	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.017	88	118208	16.751	ng	98
3) Pyridine	3.393	79	345984	17.666	ng	100
4) n-Nitrosodimethylamine	3.317	42	154442	19.362	ng	# 96
6) Aniline	6.699	93	256040	12.037	ng	99
8) 2-Chlorophenol	6.928	128	429257	22.245	ng	98
9) Benzaldehyde	6.516	77	163837	10.113	ng	97
10) Phenol	6.593	94	508661	22.014	ng	99
11) bis(2-Chloroethyl)ether	6.799	93	411525	20.552	ng	100
12) 1,3-Dichlorobenzene	7.234	146	474469	20.108	ng	99
13) 1,4-Dichlorobenzene	7.381	146	480881	20.123	ng	99
14) 1,2-Dichlorobenzene	7.687	146	467948	20.663	ng	99
15) Benzyl Alcohol	7.605	79	372737	22.893	ng	100
16) 2,2'-oxybis(1-Chloropr...	7.869	45	488602	20.350	ng	99
17) 2-Methylphenol	7.816	107	345118	21.498	ng	98
18) Hexachloroethane	8.393	117	161537	18.083	ng	98
19) n-Nitroso-di-n-propyla...	8.157	70	294609	18.249	ng	99
20) 3+4-Methylphenols	8.140	107	472164	21.154	ng	99
22) Acetophenone	8.169	105	585423	19.849	ng	99
24) Nitrobenzene	8.534	77	468483	19.888	ng	99
25) Isophorone	9.063	82	855058	19.753	ng	98
26) 2-Nitrophenol	9.234	139	235865	20.356	ng	98
27) 2,4-Dimethylphenol	9.322	122	346613	26.502	ng	99
28) bis(2-Chloroethoxy)met...	9.546	93	521338	19.667	ng	98
29) 2,4-Dichlorophenol	9.775	162	449118	22.137	ng	100
30) 1,2,4-Trichlorobenzene	9.969	180	477568	20.379	ng	98
31) Naphthalene	10.151	128	1372775	22.117	ng	100
32) Benzoic acid	9.499	122	331172	21.205	ng	99
33) 4-Chloroaniline	10.287	127	276874	11.997	ng	100
34) Hexachlorobutadiene	10.428	225	299760	20.482	ng	100
35) Caprolactam	11.093	113	136270	20.734	ng	98
36) 4-Chloro-3-methylphenol	11.434	107	434261	21.078	ng	98
37) 2-Methylnaphthalene	11.775	142	975989	21.565	ng	99
38) 1-Methylnaphthalene	11.998	142	899088	20.087	ng	100
40) 1,2,4,5-Tetrachloroben...	12.157	216	538873	20.852	ng	100
41) Hexachlorocyclopentadiene	12.128	237	14556	1.785	ng	99
43) 2,4,6-Trichlorophenol	12.422	196	379467	21.419	ng	99

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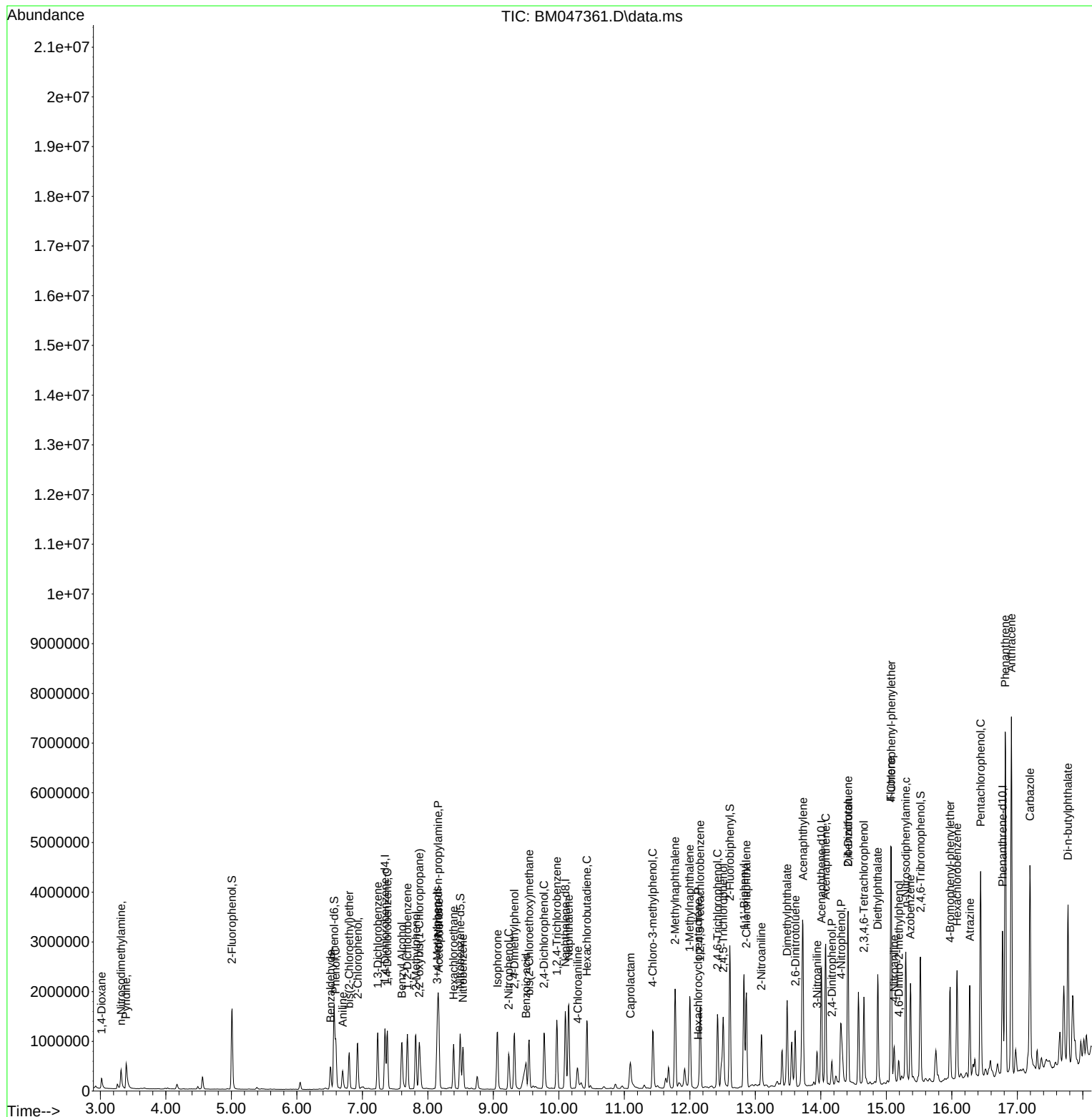
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.510	196	409487	20.813	ng	99
46) 1,1'-Biphenyl	12.828	154	1232067	20.218	ng	100
47) 2-Chloronaphthalene	12.863	162	965510	20.218	ng	99
48) 2-Nitroaniline	13.092	65	289511	21.643	ng	99
49) Acenaphthylene	13.728	152	2527867	36.421	ng	99
50) Dimethylphthalate	13.487	163	1192878	19.837	ng	100
51) 2,6-Dinitrotoluene	13.610	165	262172	18.725	ng	98
52) Acenaphthene	14.075	154	1039600	23.595	ng	99
53) 3-Nitroaniline	13.945	138	191533	14.825	ng	99
54) 2,4-Dinitrophenol	14.169	184	116257	13.483	ng	98
55) Dibenzofuran	14.416	168	1611262	22.704	ng	100
56) 4-Nitrophenol	14.310	139	460340	45.583	ng	98
57) 2,4-Dinitrotoluene	14.416	165	351309	19.383	ng	97
58) Fluorene	15.069	166	1408365	24.520	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.657	232	370489	22.926	ng	100
60) Diethylphthalate	14.869	149	1192375	20.101	ng	99
61) 4-Chlorophenyl-phenyle...	15.075	204	578512	19.800	ng	99
62) 4-Nitroaniline	15.122	138	205075	17.159	ng	98
63) Azobenzene	15.369	77	1056256	20.448	ng	95
65) 4,6-Dinitro-2-methylph...	15.192	198	89508	8.209	ng	98
66) n-Nitrosodiphenylamine	15.298	169	1030809	21.205	ng	99
67) 4-Bromophenyl-phenylether	15.975	248	383958	20.487	ng	97
68) Hexachlorobenzene	16.080	284	436137	20.148	ng	98
69) Atrazine	16.275	200	354967	22.340	ng	99
70) Pentachlorophenol	16.439	266	694721	49.640	ng	99
71) Phenanthrene	16.816	178	4168851	48.319	ng	99
72) Anthracene	16.910	178	4354622	51.775	ng	99
73) Carbazole	17.192	167	2705780	33.525	ng	99
74) Di-n-butylphthalate	17.774	149	2010302	20.455	ng	100
75) Fluoranthene	18.863	202	10671856	105.343	ng	98
78) Pyrene	19.233	202	11811096	135.526	ng	98
80) Butylbenzylphthalate	20.163	149	768314	22.529	ng	99
81) Benzo(a)anthracene	21.010	228	5558928	75.018	ng	94
82) 3,3'-Dichlorobenzidine	20.951	252	96524	3.872	ng	# 94
83) Chrysene	21.068	228	6948295	99.682	ng	97
84) Bis(2-ethylhexyl)phtha...	20.957	149	1113483	23.101	ng	99
85) Di-n-octyl phthalate	21.992	149	1755146	21.634	ng	99
87) Indeno(1,2,3-cd)pyrene	26.744	276	4088043	49.975	ng	97
88) Benzo(b)fluoranthene	22.898	252	9356072	122.722	ng	99
89) Benzo(k)fluoranthene	22.945	252	3638700	50.348	ng	98
90) Benzo(a)pyrene	23.604	252	4291868	65.098	ng	98
91) Dibenzo(a,h)anthracene	26.774	278	2009841	30.328	ng	99
92) Benzo(g,h,i)perylene	27.674	276	3377425	49.023	ng	99

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

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