

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090524\
 Data File : BM047450.D
 Acq On : 05 Sep 2024 16:45
 Operator : RC/JU
 Sample : P3836-01MS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DN-B-38(B-8)MS

Quant Time: Sep 05 17:47:53 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	7.339	152	301006	20.000	ng	-0.01
21) Naphthalene-d8	10.092	136	1078883	20.000	ng	-0.01
39) Acenaphthene-d10	14.004	164	654737	20.000	ng	0.00
64) Phenanthrene-d10	16.774	188	1136195	20.000	ng	0.00
76) Chrysene-d12	21.021	240	786808	20.000	ng	0.00
86) Perylene-d12	23.721	264	968604	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.010	112	1676960	106.886	ng	0.00
7) Phenol-d6	6.575	99	2159512	101.040	ng	0.00
23) Nitrobenzene-d5	8.492	82	1391265	71.551	ng	0.00
42) 2,4,6-Tribromophenol	15.521	330	795452	100.699	ng	0.00
45) 2-Fluorobiphenyl	12.604	172	3047505	71.517	ng	-0.01
79) Terphenyl-d14	19.439	244	2931530	66.319	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.016	88	221892	34.553	ng	99
3) Pyridine	3.387	79	611186	34.292	ng	99
4) n-Nitrosodimethylamine	3.310	42	290181	39.976	ng	97
6) Aniline	6.692	93	575668	29.739	ng	98
8) 2-Chlorophenol	6.922	128	829389	47.230	ng	97
9) Benzaldehyde	6.510	77	212186	15.497	ng	98
10) Phenol	6.598	94	937623	44.590	ng	100
11) bis(2-Chloroethyl)ether	6.792	93	793769	43.560	ng	99
12) 1,3-Dichlorobenzene	7.228	146	948043	44.149	ng	99
13) 1,4-Dichlorobenzene	7.375	146	959578	44.124	ng	99
14) 1,2-Dichlorobenzene	7.681	146	920889	44.683	ng	99
15) Benzyl Alcohol	7.604	79	713987	48.186	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.863	45	928856	42.511	ng	98
17) 2-Methylphenol	7.816	107	651128	44.569	ng	99
18) Hexachloroethane	8.381	117	361509	44.468	ng	96
19) n-Nitroso-di-n-propyla...	8.151	70	566598	38.565	ng	97
20) 3+4-Methylphenols	8.145	107	859287	42.303	ng	89
22) Acetophenone	8.163	105	1121876	44.806	ng	# 96
24) Nitrobenzene	8.534	77	877502	43.879	ng	98
25) Isophorone	9.057	82	1604315	43.655	ng	97
26) 2-Nitrophenol	9.233	139	486143	49.421	ng	99
27) 2,4-Dimethylphenol	9.322	122	646893	58.261	ng	99
28) bis(2-Chloroethoxy)met...	9.539	93	1014090	45.062	ng	98
29) 2,4-Dichlorophenol	9.786	162	877938	50.972	ng	99
30) 1,2,4-Trichlorobenzene	9.963	180	928634	46.678	ng	99
31) Naphthalene	10.145	128	2400102	45.547	ng	99
32) Benzoic acid	9.557	122	507113	38.248	ng	96
33) 4-Chloroaniline	10.292	127	506515	25.852	ng	99
34) Hexachlorobutadiene	10.416	225	627754	50.525	ng	99
35) Caprolactam	11.116	113	232653	41.697	ng	97
36) 4-Chloro-3-methylphenol	11.451	107	759161	43.403	ng	98
37) 2-Methylnaphthalene	11.769	142	1707296	44.436	ng	99
38) 1-Methylnaphthalene	11.998	142	1596804	42.022	ng	99
40) 1,2,4,5-Tetrachloroben...	12.157	216	1028726	53.060	ng	99
41) Hexachlorocyclopentadiene	12.122	237	574118	93.826	ng	99
43) 2,4,6-Trichlorophenol	12.427	196	685162	51.552	ng	99

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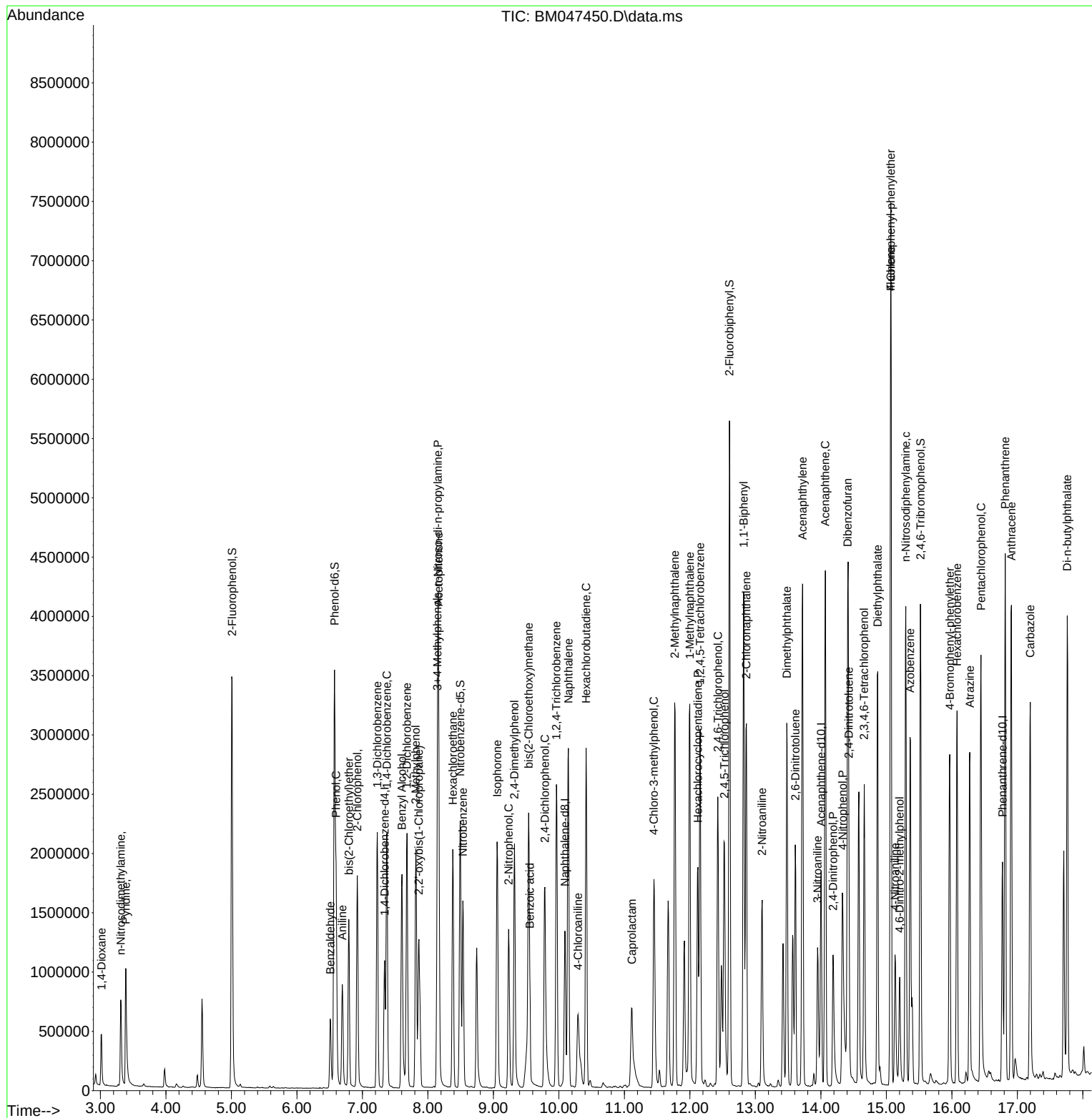
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.527	196	707198	47.912	ng	97
46) 1,1'-Biphenyl	12.821	154	2201374	48.152	ng	100
47) 2-Chloronaphthalene	12.863	162	1683666	46.995	ng	99
48) 2-Nitroaniline	13.104	65	461326	45.969	ng	93
49) Acenaphthylene	13.721	152	2575045	49.454	ng	99
50) Dimethylphthalate	13.480	163	2162653	47.937	ng	100
51) 2,6-Dinitrotoluene	13.610	165	469890	44.735	ng	98
52) Acenaphthene	14.068	154	1559197	47.170	ng	99
53) 3-Nitroaniline	13.951	138	372429	38.424	ng	96
54) 2,4-Dinitrophenol	14.186	184	383515	59.288	ng	95
55) Dibenzofuran	14.410	168	2477498	46.533	ng	98
56) 4-Nitrophenol	14.333	139	526260	69.460	ng	98
57) 2,4-Dinitrotoluene	14.427	165	584893	43.015	ng	# 83
58) Fluorene	15.068	166	1890463	43.873	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.663	232	565644	46.656	ng	97
60) Diethylphthalate	14.868	149	2024466	45.492	ng	99
61) 4-Chlorophenyl-phenyle...	15.068	204	1024034	46.717	ng	95
62) 4-Nitroaniline	15.133	138	360728	40.232	ng	97
63) Azobenzene	15.363	77	1830269	47.230	ng	98
65) 4,6-Dinitro-2-methylph...	15.204	198	251383	36.048	ng	100
66) n-Nitrosodiphenylamine	15.298	169	1658173	53.336	ng	100
67) 4-Bromophenyl-phenylether	15.968	248	624445	52.096	ng	94
68) Hexachlorobenzene	16.080	284	682505	49.299	ng	96
69) Atrazine	16.274	200	588077	57.870	ng	99
70) Pentachlorophenol	16.445	266	777120	86.823	ng	98
71) Phenanthrene	16.815	178	2687270	48.701	ng	100
72) Anthracene	16.909	178	2693429	50.073	ng	100
73) Carbazole	17.198	167	2326184	45.066	ng	100
74) Di-n-butylphthalate	17.762	149	2959801	47.090	ng	100
75) Fluoranthene	18.856	202	2816439	43.470	ng	99
77) Benzidine	19.068	184	1305156	99.375	ng	100
78) Pyrene	19.221	202	2861543	47.419	ng	99
80) Butylbenzylphthalate	20.150	149	1114166	47.181	ng	96
81) Benzo(a)anthracene	21.003	228	2518100	49.076	ng	99
82) 3,3'-Dichlorobenzidine	20.944	252	814179	47.171	ng	99
83) Chrysene	21.056	228	2356497	48.823	ng	99
84) Bis(2-ethylhexyl)phtha...	20.944	149	1545715	46.312	ng	# 97
85) Di-n-octyl phthalate	21.968	149	2639767	46.991	ng	98
87) Indeno(1,2,3-cd)pyrene	26.744	276	3186469	52.552	ng	98
88) Benzo(b)fluoranthene	22.880	252	2589292	45.820	ng	100
89) Benzo(k)fluoranthene	22.933	252	2531824	47.262	ng	99
90) Benzo(a)pyrene	23.597	252	2504708	51.253	ng	99
91) Dibenzo(a,h)anthracene	26.774	278	2529785	51.501	ng	100
92) Benzo(g,h,i)perylene	27.685	276	2465224	48.274	ng	98

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

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