

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090424\
 Data File : BM047441.D
 Acq On : 04 Sep 2024 16:43
 Operator : RC/JU
 Sample : P3825-01MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DN-W-01(B-1)MSD

Quant Time: Sep 04 17:28:35 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	7.340	152	211500	20.000	ng	-0.01
21) Naphthalene-d8	10.098	136	752008	20.000	ng	0.00
39) Acenaphthene-d10	14.004	164	485384	20.000	ng	0.00
64) Phenanthrene-d10	16.774	188	945169	20.000	ng	0.00
76) Chrysene-d12	21.021	240	805183	20.000	ng	0.00
86) Perylene-d12	23.727	264	947731	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.010	112	1402347	127.210	ng	0.00
7) Phenol-d6	6.581	99	1723742	114.782	ng	0.01
23) Nitrobenzene-d5	8.492	82	1119883	82.629	ng	0.00
42) 2,4,6-Tribromophenol	15.527	330	724064	123.643	ng	0.00
45) 2-Fluorobiphenyl	12.610	172	2774848	87.839	ng	0.00
79) Terphenyl-d14	19.439	244	3560325	78.705	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.016	88	218766	48.483	ng	97
3) Pyridine	3.393	79	571405	45.627	ng	99
4) n-Nitrosodimethylamine	3.310	42	259507	50.880	ng	98
6) Aniline	6.704	93	296379	21.790	ng	99
8) 2-Chlorophenol	6.928	128	685887	55.588	ng	98
9) Benzaldehyde	6.510	77	181212	19.650	ng	96
10) Phenol	6.604	94	796832	53.931	ng	99
11) bis(2-Chloroethyl)ether	6.793	93	768283	60.003	ng	95
12) 1,3-Dichlorobenzene	7.228	146	818427	54.242	ng	98
13) 1,4-Dichlorobenzene	7.375	146	839910	54.966	ng	99
14) 1,2-Dichlorobenzene	7.681	146	792836	54.749	ng	100
15) Benzyl Alcohol	7.604	79	581018	55.806	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.863	45	766486	49.925	ng	98
17) 2-Methylphenol	7.822	107	534094	52.029	ng	99
18) Hexachloroethane	8.387	117	281946	49.358	ng	94
19) n-Nitroso-di-n-propyla...	8.157	70	478736	46.375	ng	98
20) 3+4-Methylphenols	8.157	107	714744	50.078	ng	# 62
22) Acetophenone	8.169	105	962917	55.173	ng	# 97
24) Nitrobenzene	8.540	77	721022	51.726	ng	96
25) Isophorone	9.063	82	1315633	51.361	ng	98
26) 2-Nitrophenol	9.240	139	381071	55.579	ng	95
27) 2,4-Dimethylphenol	9.328	122	527512	68.160	ng	100
28) bis(2-Chloroethoxy)met...	9.540	93	805103	51.326	ng	98
29) 2,4-Dichlorophenol	9.798	162	693996	57.806	ng	97
30) 1,2,4-Trichlorobenzene	9.969	180	778336	56.129	ng	99
31) Naphthalene	10.145	128	2054775	55.943	ng	99
32) Benzoic acid	9.551	122	443917	48.035	ng	97
33) 4-Chloroaniline	10.304	127	419065	30.685	ng	98
34) Hexachlorobutadiene	10.422	225	519576	59.996	ng	99
35) Caprolactam	11.116	113	189830	48.810	ng	99
36) 4-Chloro-3-methylphenol	11.457	107	611837	50.185	ng	98
37) 2-Methylnaphthalene	11.775	142	1451811	54.211	ng	99
38) 1-Methylnaphthalene	11.998	142	1353363	51.096	ng	99
40) 1,2,4,5-Tetrachloroben...	12.157	216	864321	60.135	ng	99
41) Hexachlorocyclopentadiene	12.122	237	327969	72.300	ng	99
43) 2,4,6-Trichlorophenol	12.433	196	574136	58.270	ng	99

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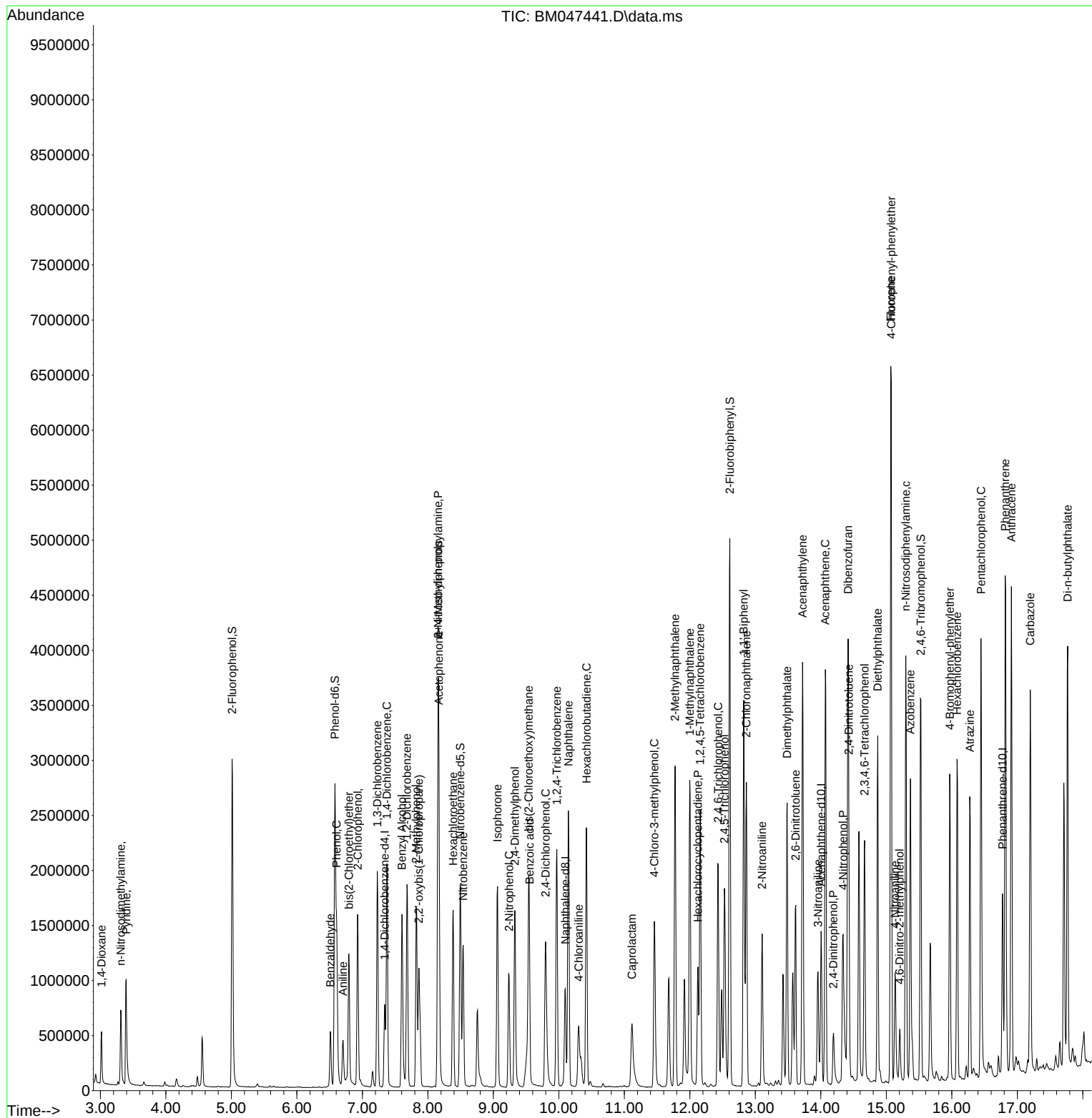
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.528	196	607125	55.483	ng	97
46) 1,1'-Biphenyl	12.822	154	1888728	55.728	ng	100
47) 2-Chloronaphthalene	12.863	162	1475855	55.567	ng	99
48) 2-Nitroaniline	13.104	65	385337	51.794	ng	95
49) Acenaphthylene	13.722	152	2325826	60.252	ng	99
50) Dimethylphthalate	13.486	163	1784182	53.347	ng	99
51) 2,6-Dinitrotoluene	13.616	165	390714	50.175	ng	93
52) Acenaphthene	14.069	154	1378957	56.273	ng	98
53) 3-Nitroaniline	13.957	138	328474	45.713	ng	94
54) 2,4-Dinitrophenol	14.192	184	163210	34.034	ng	93
55) Dibenzofuran	14.416	168	2229426	56.483	ng	99
56) 4-Nitrophenol	14.339	139	540181	96.174	ng	96
57) 2,4-Dinitrotoluene	14.427	165	524630	52.045	ng	# 84
58) Fluorene	15.069	166	1737227	54.383	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.669	232	517585	57.588	ng	95
60) Diethylphthalate	14.869	149	1704168	51.656	ng	98
61) 4-Chlorophenyl-phenyle...	15.074	204	916339	56.389	ng	95
62) 4-Nitroaniline	15.133	138	323139	48.614	ng	96
63) Azobenzene	15.369	77	1530397	53.271	ng	96
65) 4,6-Dinitro-2-methylph...	15.204	198	130385	22.476	ng	99
66) n-Nitrosodiphenylamine	15.298	169	1483954	57.379	ng	99
67) 4-Bromophenyl-phenylether	15.968	248	585432	58.713	ng	95
68) Hexachlorobenzene	16.080	284	643563	55.882	ng	96
69) Atrazine	16.274	200	528639	62.535	ng	99
70) Pentachlorophenol	16.445	266	808873	108.635	ng	99
71) Phenanthrene	16.816	178	2852169	62.137	ng	100
72) Anthracene	16.910	178	2750984	61.479	ng	100
73) Carbazole	17.198	167	2377629	55.372	ng	100
74) Di-n-butylphthalate	17.768	149	2815960	53.856	ng	100
75) Fluoranthene	18.857	202	3634321	67.431	ng	99
77) Benzidine	19.074	184	66097	4.918	ng	# 87
78) Pyrene	19.221	202	3731957	60.432	ng	100
80) Butylbenzylphthalate	20.156	149	1248790	51.675	ng	95
81) Benzo(a)anthracene	21.009	228	3238193	61.670	ng	99
82) 3,3'-Dichlorobenzidine	20.945	252	717775	40.637	ng	99
83) Chrysene	21.062	228	3083553	62.429	ng	100
84) Bis(2-ethylhexyl)phtha...	20.951	149	2137361	62.577	ng	# 98
85) Di-n-octyl phthalate	21.980	149	3003649	52.248	ng	99
87) Indeno(1,2,3-cd)pyrene	26.768	276	3946977	66.528	ng	97
88) Benzo(b)fluoranthene	22.886	252	3402018	61.528	ng	100
89) Benzo(k)fluoranthene	22.939	252	3211879	61.277	ng	100
90) Benzo(a)pyrene	23.609	252	3198336	66.888	ng	99
91) Dibenzo(a,h)anthracene	26.791	278	3027560	62.992	ng	99
92) Benzo(g,h,i)perylene	27.697	276	3090887	61.859	ng	98

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

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