

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM090324\
 Data File : BM047428.D
 Acq On : 03 Sep 2024 16:43
 Operator : RC/JU
 Sample : P3813-01MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TR-04-083024MSD

Quant Time: Sep 03 17:20:48 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	246733	20.000	ng	-0.01
21) Naphthalene-d8	10.098	136	926995	20.000	ng	0.00
39) Acenaphthene-d10	14.010	164	574254	20.000	ng	0.00
64) Phenanthrene-d10	16.774	188	1028165	20.000	ng	0.00
76) Chrysene-d12	21.021	240	804689	20.000	ng	0.00
86) Perylene-d12	23.727	264	1000348	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.010	112	1619460	125.927	ng	0.00
7) Phenol-d6	6.575	99	2091548	119.386	ng	0.00
23) Nitrobenzene-d5	8.498	82	1389642	83.178	ng	0.00
42) 2,4,6-Tribromophenol	15.527	330	826456	119.287	ng	0.00
45) 2-Fluorobiphenyl	12.610	172	3314177	88.676	ng	0.00
79) Terphenyl-d14	19.445	244	3704259	81.937	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.011	88	210235	39.939	ng	99
3) Pyridine	3.387	79	572018	39.154	ng	99
4) n-Nitrosodimethylamine	3.310	42	270205	45.412	ng	100
6) Aniline	6.699	93	512804	32.318	ng	99
8) 2-Chlorophenol	6.928	128	770158	53.504	ng	97
9) Benzaldehyde	6.510	77	207269	19.181	ng	99
10) Phenol	6.598	94	867692	50.341	ng	99
11) bis(2-Chloroethyl)ether	6.793	93	719664	48.180	ng	100
12) 1,3-Dichlorobenzene	7.228	146	880235	50.008	ng	98
13) 1,4-Dichlorobenzene	7.375	146	895168	50.216	ng	99
14) 1,2-Dichlorobenzene	7.687	146	860481	50.936	ng	99
15) Benzyl Alcohol	7.604	79	658789	54.240	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.863	45	865301	48.314	ng	98
17) 2-Methylphenol	7.816	107	613140	51.200	ng	99
18) Hexachloroethane	8.387	117	315679	47.372	ng	97
19) n-Nitroso-di-n-propyla...	8.157	70	519558	43.143	ng	97
20) 3+4-Methylphenols	8.145	107	813767	48.874	ng	96
22) Acetophenone	8.169	105	1041208	48.397	ng	100
24) Nitrobenzene	8.540	77	820603	47.757	ng	97
25) Isophorone	9.063	82	1504648	47.651	ng	98
26) 2-Nitrophenol	9.239	139	452978	53.595	ng	96
27) 2,4-Dimethylphenol	9.322	122	613843	64.343	ng	98
28) bis(2-Chloroethoxy)met...	9.539	93	938424	48.532	ng	98
29) 2,4-Dichlorophenol	9.781	162	824971	55.745	ng	99
30) 1,2,4-Trichlorobenzene	9.969	180	878574	51.398	ng	98
31) Naphthalene	10.145	128	2289676	50.571	ng	99
32) Benzoic acid	9.551	122	501816	44.050	ng	95
33) 4-Chloroaniline	10.287	127	449220	26.684	ng	99
34) Hexachlorobutadiene	10.422	225	580725	54.398	ng	98
35) Caprolactam	11.110	113	239344	49.925	ng	100
36) 4-Chloro-3-methylphenol	11.451	107	744549	49.542	ng	99
37) 2-Methylnaphthalene	11.775	142	1645861	49.856	ng	99
38) 1-Methylnaphthalene	12.004	142	1525523	46.724	ng	100
40) 1,2,4,5-Tetrachloroben...	12.163	216	985131	57.933	ng	99
41) Hexachlorocyclopentadiene	12.122	237	124686	23.233	ng	99
43) 2,4,6-Trichlorophenol	12.428	196	665617	57.100	ng	99

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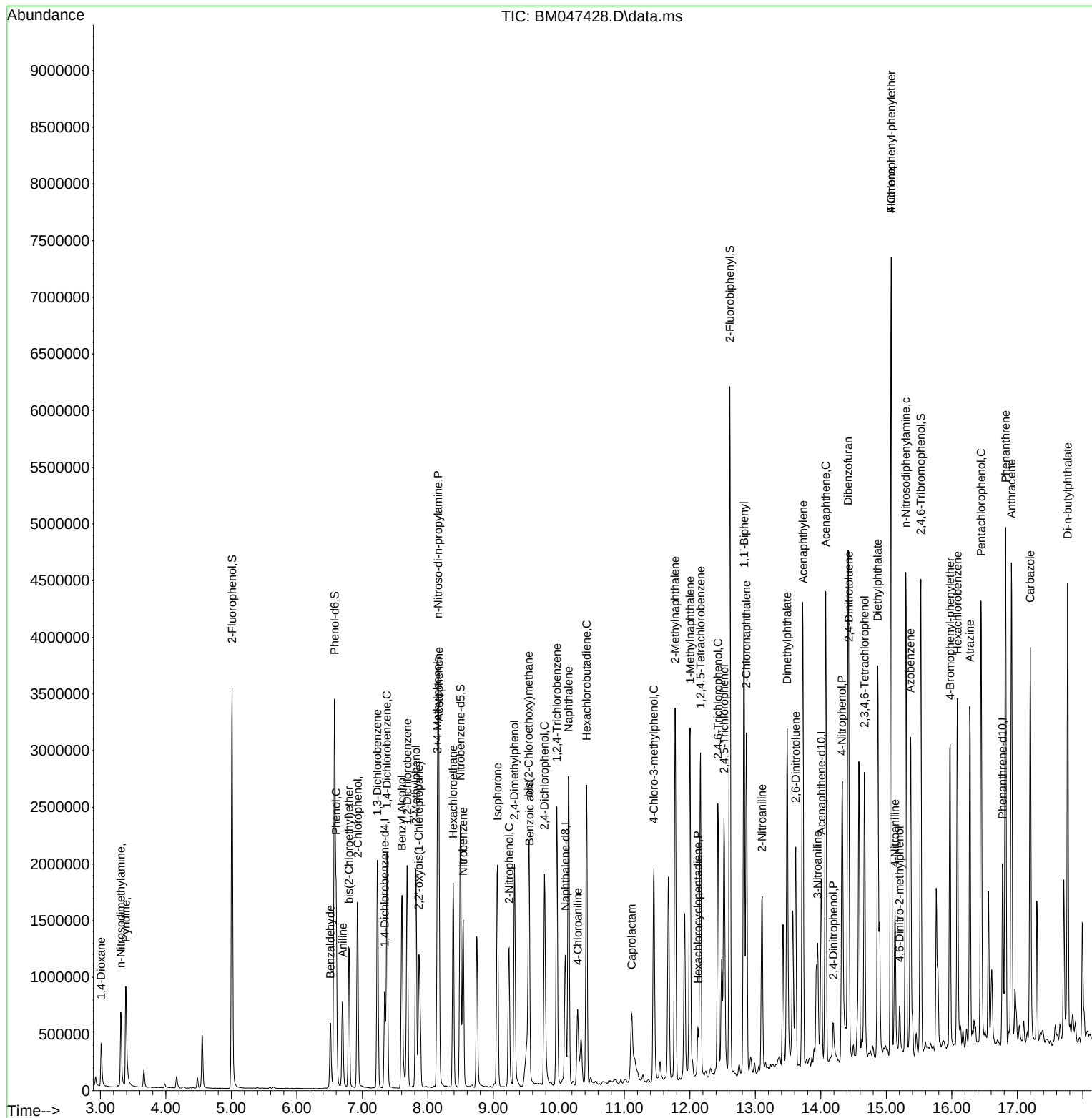
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.522	196	704783	54.441	ng	98
46) 1,1'-Biphenyl	12.828	154	2098129	52.326	ng	100
47) 2-Chloronaphthalene	12.863	162	1616917	51.457	ng	99
48) 2-Nitroaniline	13.104	65	462269	52.519	ng	94
49) Acenaphthylene	13.727	152	2560758	56.072	ng	99
50) Dimethylphthalate	13.486	163	2046392	51.718	ng	100
51) 2,6-Dinitrotoluene	13.616	165	454911	49.379	ng	96
52) Acenaphthene	14.074	154	1525477	52.618	ng	99
53) 3-Nitroaniline	13.951	138	317056	37.296	ng	97
54) 2,4-Dinitrophenol	14.192	184	120267	21.198	ng	91
55) Dibenzofuran	14.416	168	2383874	51.050	ng	99
56) 4-Nitrophenol	14.322	139	637589	95.949	ng	99
57) 2,4-Dinitrotoluene	14.427	165	578523	48.509	ng	# 85
58) Fluorene	15.074	166	1851148	48.981	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.669	232	568984	53.509	ng	98
60) Diethylphthalate	14.869	149	1938208	49.658	ng	99
61) 4-Chlorophenyl-phenyle...	15.074	204	967456	50.322	ng	97
62) 4-Nitroaniline	15.133	138	367897	46.782	ng	97
63) Azobenzene	15.369	77	1649946	48.544	ng	98
65) 4,6-Dinitro-2-methylph...	15.204	198	101262	16.046	ng	93
66) n-Nitrosodiphenylamine	15.298	169	1612056	57.300	ng	98
67) 4-Bromophenyl-phenylether	15.974	248	607685	56.025	ng	96
68) Hexachlorobenzene	16.086	284	666772	53.223	ng	98
69) Atrazine	16.274	200	597200	64.942	ng	99
70) Pentachlorophenol	16.445	266	816722	100.835	ng	99
71) Phenanthrene	16.821	178	2833355	56.744	ng	99
72) Anthracene	16.910	178	2783061	57.176	ng	100
73) Carbazole	17.198	167	2405748	51.504	ng	100
74) Di-n-butylphthalate	17.768	149	3007028	52.868	ng	100
75) Fluoranthene	18.857	202	3359959	57.308	ng	99
77) Benzidine	19.068	184	1895250	141.098	ng	99
78) Pyrene	19.227	202	3475545	56.314	ng	99
80) Butylbenzylphthalate	20.156	149	1192933	49.394	ng	98
81) Benzo(a)anthracene	21.009	228	3056897	58.253	ng	98
82) 3,3'-Dichlorobenzidine	20.945	252	996392	56.445	ng	98
83) Chrysene	21.062	228	2889531	58.536	ng	99
84) Bis(2-ethylhexyl)phtha...	20.951	149	1636320	47.937	ng	# 99
85) Di-n-octyl phthalate	21.980	149	2981200	51.889	ng	98
87) Indeno(1,2,3-cd)pyrene	26.762	276	3924742	62.674	ng	97
88) Benzo(b)fluoranthene	22.886	252	3350683	57.412	ng	100
89) Benzo(k)fluoranthene	22.945	252	3009422	54.395	ng	99
90) Benzo(a)pyrene	23.609	252	3187851	63.162	ng	99
91) Dibenzo(a,h)anthracene	26.797	278	2962448	58.395	ng	99
92) Benzo(g,h,i)perylene	27.697	276	3118793	59.135	ng	98

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

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