

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040425\
 Data File : BM049831.D
 Acq On : 04 Apr 2025 19:56
 Operator : RC/JU
 Sample : Q1694-01MS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 TP-18MS

Quant Time: Apr 04 22:57:18 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM031325.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 13 14:55:34 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.792	152	356782	20.000	ng	-0.01
21) Naphthalene-d8	10.586	136	1294303	20.000	ng	-0.02
39) Acenaphthene-d10	14.433	164	862231	20.000	ng	-0.01
64) Phenanthrene-d10	17.174	188	1735840	20.000	ng	-0.01
76) Chrysene-d12	21.415	240	1721481	20.000	ng	-0.01
86) Perylene-d12	24.421	264	1866539	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	2131663	100.813	ng	-0.01
7) Phenol-d6	6.969	99	2740294	103.268	ng	-0.01
23) Nitrobenzene-d5	8.945	82	1588023	62.993	ng	-0.02
42) 2,4,6-Tribromophenol	15.921	330	1411590	140.202	ng	-0.01
45) 2-Fluorobiphenyl	13.057	172	4344581	63.104	ng	-0.01
79) Terphenyl-d14	19.798	244	7051205	69.023	ng	-0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.269	88	327074	35.732	ng	96
3) Pyridine	3.669	79	905148	38.720	ng	98
4) n-Nitrosodimethylamine	3.575	42	369701	35.742	ng	89
6) Aniline	7.116	93	1180028	53.080	ng	97
8) 2-Chlorophenol	7.357	128	1045795	49.944	ng	97
9) Benzaldehyde	6.928	77	602975	45.965	ng	95
10) Phenol	6.992	94	1255026	48.777	ng	95
11) bis(2-Chloroethyl)ether	7.210	93	928576	44.566	ng	97
12) 1,3-Dichlorobenzene	7.681	146	1191597	46.827	ng	97
13) 1,4-Dichlorobenzene	7.828	146	1200257	46.754	ng	98
14) 1,2-Dichlorobenzene	8.139	146	1164321	47.723	ng	99
15) Benzyl Alcohol	8.028	79	844425	50.709	ng	93
16) 2,2'-oxybis(1-Chloropr...	8.316	45	1096453	44.774	ng	98
17) 2-Methylphenol	8.239	107	813160	53.232	ng	96
18) Hexachloroethane	8.869	117	379254	41.834	ng	89
19) n-Nitroso-di-n-propyla...	8.598	70	710167	46.154	ng	96
20) 3+4-Methylphenols	8.563	107	1111107	54.276	ng	95
22) Acetophenone	8.610	105	1422024	42.610	ng	97
24) Nitrobenzene	8.986	77	1020526	42.932	ng	95
25) Isophorone	9.516	82	1950869	50.402	ng	98
26) 2-Nitrophenol	9.698	139	540494	53.422	ng	96
27) 2,4-Dimethylphenol	9.769	122	887210	68.892	ng	96
28) bis(2-Chloroethoxy)met...	9.998	93	1219253	44.725	ng	98
29) 2,4-Dichlorophenol	10.239	162	1083945	51.654	ng	97
30) 1,2,4-Trichlorobenzene	10.451	180	1231870	46.920	ng	99
31) Naphthalene	10.639	128	3041164	45.695	ng	99
32) Benzoic acid	9.916	122	699773	63.892	ng	96
33) 4-Chloroaniline	10.745	127	856792	37.722	ng	97
34) Hexachlorobutadiene	10.928	225	764172	47.148	ng	99
35) Caprolactam	11.528	113	296534	60.865	ng	# 87
36) 4-Chloro-3-methylphenol	11.880	107	950287	52.769	ng	97
37) 2-Methylnaphthalene	12.251	142	2045877	43.962	ng	98
38) 1-Methylnaphthalene	12.475	142	2129077	47.419	ng	98
40) 1,2,4,5-Tetrachloroben...	12.622	216	1428498	45.104	ng	98
41) Hexachlorocyclopentadiene	12.604	237	575323	49.034	ng	98
43) 2,4,6-Trichlorophenol	12.863	196	937302	52.841	ng	97

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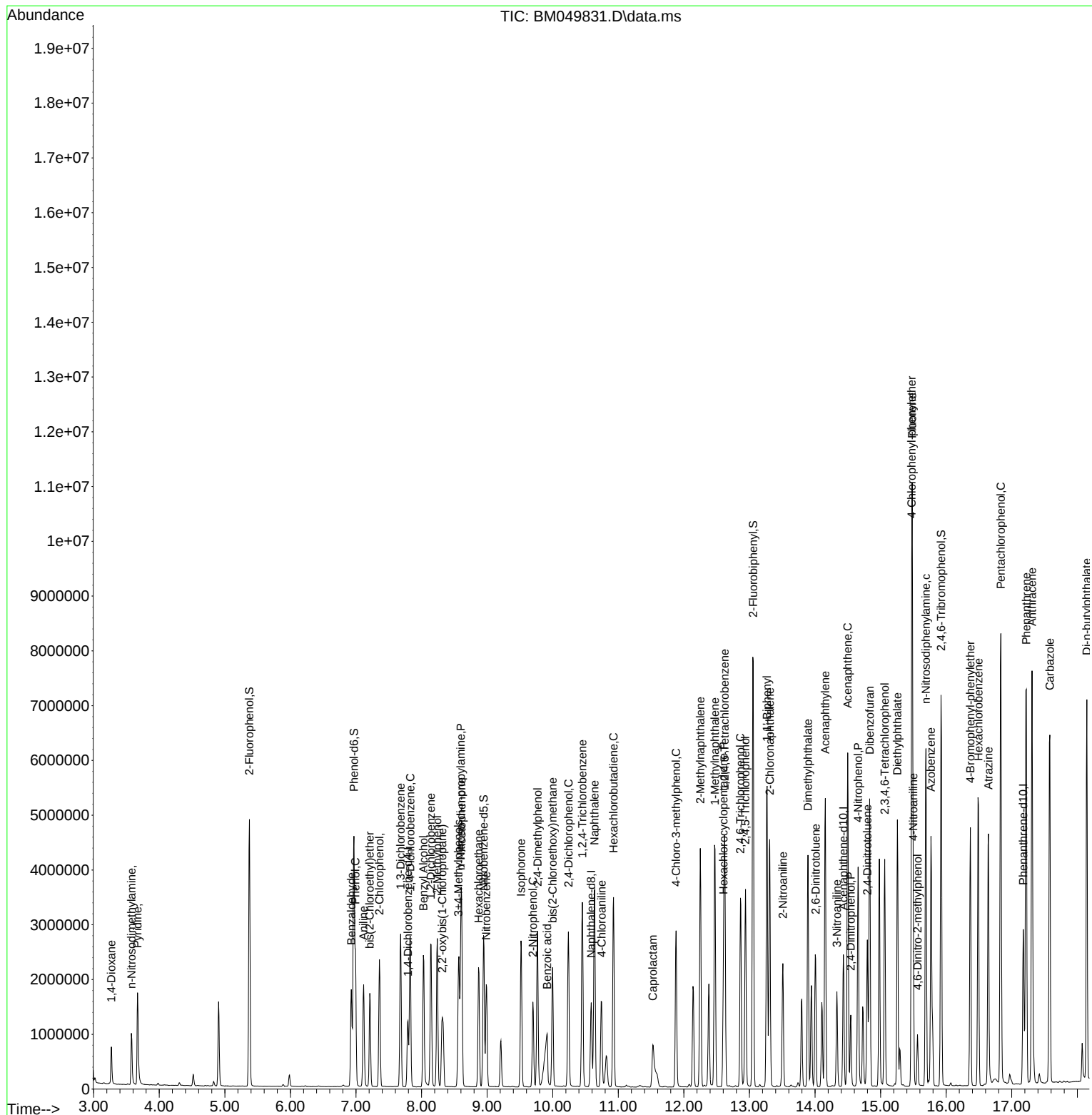
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.939	196	1020135	52.702	ng	97
46) 1,1'-Biphenyl	13.269	154	2987687	44.959	ng	98
47) 2-Chloronaphthalene	13.310	162	2383788	46.086	ng	100
48) 2-Nitroaniline	13.510	65	577604	49.744	ng	88
49) Acenaphthylene	14.157	152	3769193	52.133	ng	99
50) Dimethylphthalate	13.892	163	2800425	47.174	ng	100
51) 2,6-Dinitrotoluene	14.010	165	601714	50.586	ng	# 86
52) Acenaphthene	14.498	154	2182044	46.011	ng	98
53) 3-Nitroaniline	14.333	138	447415	39.527	ng	98
54) 2,4-Dinitrophenol	14.545	184	308689	49.273	ng	91
55) Dibenzofuran	14.833	168	3584928	45.365	ng	97
56) 4-Nitrophenol	14.657	139	1123672	114.046	ng	88
57) 2,4-Dinitrotoluene	14.798	165	850613	55.574	ng	92
58) Fluorene	15.480	166	3001109	46.645	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.063	232	871198	53.768	ng	97
60) Diethylphthalate	15.257	149	2672859	47.806	ng	97
61) 4-Chlorophenyl-phenyle...	15.474	204	1639431	46.663	ng	98
62) 4-Nitroaniline	15.498	138	572410	42.933	ng	91
63) Azobenzene	15.768	77	2601003	48.704	ng	94
65) 4,6-Dinitro-2-methylph...	15.563	198	226523	27.920	ng	97
66) n-Nitrosodiphenylamine	15.692	169	2516077	47.841	ng	98
67) 4-Bromophenyl-phenylether	16.368	248	999084	49.851	ng	97
68) Hexachlorobenzene	16.492	284	1133078	50.276	ng	95
69) Atrazine	16.645	200	925169	79.821	ng	99
70) Pentachlorophenol	16.833	266	1588260	112.143	ng	98
71) Phenanthrene	17.221	178	4636151	48.099	ng	100
72) Anthracene	17.309	178	4759687	50.990	ng	100
73) Carbazole	17.580	167	4282604	50.659	ng	99
74) Di-n-butylphthalate	18.145	149	4962054	53.553	ng	99
75) Fluoranthene	19.233	202	5564095	50.141	ng	99
77) Benzidine	19.421	184	5144221	248.731	ng	99
78) Pyrene	19.598	202	5692052	47.326	ng	100
80) Butylbenzylphthalate	20.492	149	2186734	58.586	ng	97
81) Benzo(a)anthracene	21.397	228	5694517	49.525	ng	99
82) 3,3'-Dichlorobenzidine	21.315	252	2248387	60.060	ng	98
83) Chrysene	21.456	228	5234523	48.024	ng	100
84) Bis(2-ethylhexyl)phtha...	21.321	149	3186763	54.784	ng	# 95
85) Di-n-octyl phthalate	22.456	149	5683539	62.863	ng	96
87) Indeno(1,2,3-cd)pyrene	27.826	276	6754576	50.278	ng	99
88) Benzo(b)fluoranthene	23.474	252	5602460	44.564	ng	99
89) Benzo(k)fluoranthene	23.538	252	5581332	45.118	ng	99
90) Benzo(a)pyrene	24.285	252	5372649	50.878	ng	99
91) Dibenzo(a,h)anthracene	27.879	278	5572457	49.785	ng	99
92) Benzo(g,h,i)perylene	28.885	276	5283895	46.866	ng	100

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

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