

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020525\
 Data File : BM049570.D
 Acq On : 05 Feb 2025 14:58
 Operator : RC/JU
 Sample : SSTDICC060
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC060

Quant Time: Feb 06 04:41:45 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM020525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 06 04:35:49 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	261635	20.000	ng	0.00
21) Naphthalene-d8	10.628	136	971391	20.000	ng	0.00
39) Acenaphthene-d10	14.469	164	618462	20.000	ng	0.00
64) Phenanthrene-d10	17.210	188	1164611	20.000	ng	0.00
76) Chrysene-d12	21.445	240	1132403	20.000	ng	0.00
86) Perylene-d12	24.480	264	926026	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	2038411	125.287	ng	0.00
7) Phenol-d6	6.987	99	2719270	127.572	ng	0.00
23) Nitrobenzene-d5	8.981	82	2434266	128.900	ng	0.00
42) 2,4,6-Tribromophenol	15.951	330	1086269	148.230	ng	0.00
45) 2-Fluorobiphenyl	13.098	172	6338303	132.764	ng	0.00
79) Terphenyl-d14	19.833	244	10117181	136.430	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.293	88	417562	61.040	ng	100
3) Pyridine	3.687	79	1198421	63.042	ng	98
4) n-Nitrosodimethylamine	3.599	42	553563	62.239	ng	100
6) Aniline	7.151	93	1257983	62.269	ng	99
8) 2-Chlorophenol	7.393	128	1020217	62.798	ng	99
9) Benzaldehyde	6.963	77	596122	52.818	ng	97
10) Phenol	7.010	94	1322610	62.997	ng	100
11) bis(2-Chloroethyl)ether	7.251	93	1049352	61.780	ng	100
12) 1,3-Dichlorobenzene	7.716	146	1158517	61.436	ng	98
13) 1,4-Dichlorobenzene	7.863	146	1172598	61.477	ng	100
14) 1,2-Dichlorobenzene	8.181	146	1134645	61.608	ng	99
15) Benzyl Alcohol	8.063	79	946556	63.829	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.357	45	1397669	61.127	ng	100
17) 2-Methylphenol	8.263	107	805349	62.734	ng	99
18) Hexachloroethane	8.916	117	432749	62.113	ng	98
19) n-Nitroso-di-n-propyla...	8.639	70	889345	63.714	ng	98
20) 3+4-Methylphenols	8.592	107	1090230	64.333	ng	100
22) Acetophenone	8.645	105	1638735	62.634	ng	# 99
24) Nitrobenzene	9.022	77	1157718	63.098	ng	99
25) Isophorone	9.551	82	2133576	62.645	ng	100
26) 2-Nitrophenol	9.734	139	408973	60.890	ng	99
27) 2,4-Dimethylphenol	9.798	122	657612	61.065	ng	99
28) bis(2-Chloroethoxy)met...	10.039	93	1363833	62.174	ng	98
29) 2,4-Dichlorophenol	10.269	162	972997	65.608	ng	99
30) 1,2,4-Trichlorobenzene	10.486	180	1140312	62.759	ng	100
31) Naphthalene	10.675	128	3222711	62.905	ng	99
32) Benzoic acid	9.934	122	449135	60.088	ng	99
33) 4-Chloroaniline	10.775	127	1201211	62.641	ng	99
34) Hexachlorobutadiene	10.975	225	696721	64.032	ng	99
35) Caprolactam	11.551	113	292307	64.607	ng	99
36) 4-Chloro-3-methylphenol	11.898	107	963434	65.141	ng	99
37) 2-Methylnaphthalene	12.292	142	2300913	63.789	ng	99
38) 1-Methylnaphthalene	12.510	142	2241932	63.881	ng	99
40) 1,2,4,5-Tetrachloroben...	12.657	216	1334952	65.651	ng	99
41) Hexachlorocyclopentadiene	12.651	237	363379	70.002	ng	98
43) 2,4,6-Trichlorophenol	12.892	196	754692	68.015	ng	98

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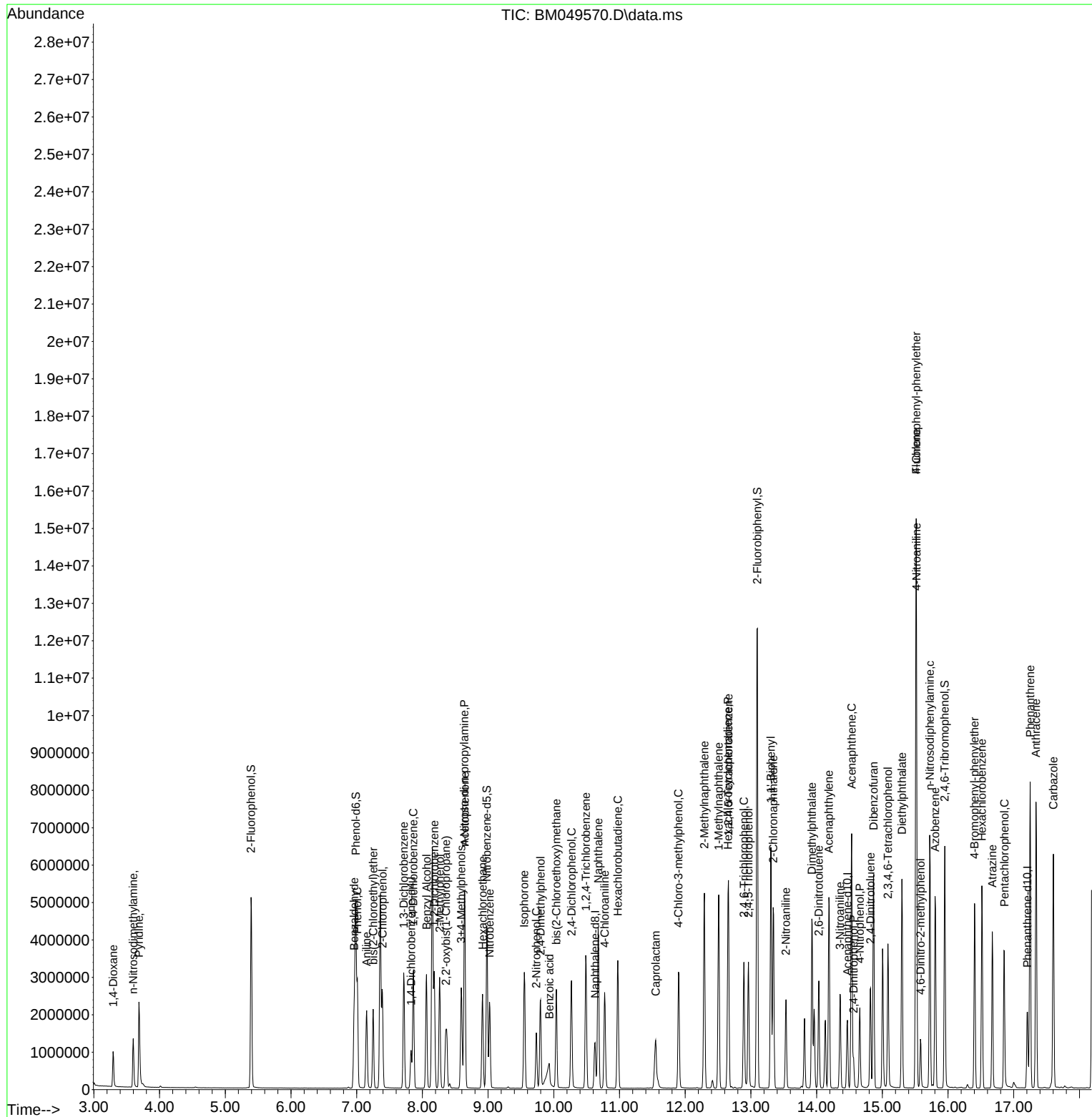
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.963	196	828431	67.938	ng	99
46) 1,1'-Biphenyl	13.304	154	3011389	63.828	ng	99
47) 2-Chloronaphthalene	13.339	162	2329169	63.808	ng	99
48) 2-Nitroaniline	13.533	65	589664	60.127	ng	99
49) Acenaphthylene	14.186	152	3444441	63.992	ng	99
50) Dimethylphthalate	13.927	163	2817395	63.434	ng	99
51) 2,6-Dinitrotoluene	14.033	165	533046	69.358	ng	99
52) Acenaphthene	14.533	154	2335450	65.368	ng	99
53) 3-Nitroaniline	14.357	138	543112	69.409	ng	100
54) 2,4-Dinitrophenol	14.563	184	140292	60.924	ng	# 78
55) Dibenzofuran	14.868	168	3615545	63.554	ng	99
56) 4-Nitrophenol	14.657	139	452050	68.533	ng	96
57) 2,4-Dinitrotoluene	14.821	165	695289	60.725	ng	93
58) Fluorene	15.515	166	3265838	67.106	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.086	232	717449	69.641	ng	98
60) Diethylphthalate	15.298	149	2854554	63.951	ng	100
61) 4-Chlorophenyl-phenyle...	15.515	204	1776049	68.566	ng	98
62) 4-Nitroaniline	15.521	138	613586	59.492	ng	97
63) Azobenzene	15.804	77	2905105	62.354	ng	99
65) 4,6-Dinitro-2-methylph...	15.580	198	235707	61.341	ng	97
66) n-Nitrosodiphenylamine	15.721	169	2408929	65.588	ng	99
67) 4-Bromophenyl-phenylether	16.404	248	897999	66.372	ng	99
68) Hexachlorobenzene	16.515	284	1019310	65.639	ng	99
69) Atrazine	16.674	200	704921	64.565	ng	96
70) Pentachlorophenol	16.857	266	602375	59.645	ng	99
71) Phenanthrene	17.251	178	4423420	66.671	ng	100
72) Anthracene	17.339	178	4372525	66.559	ng	100
73) Carbazole	17.604	167	3665434	63.570	ng	100
74) Di-n-butylphthalate	18.192	149	4884028	66.755	ng	100
75) Fluoranthene	19.262	202	5356301	68.956	ng	99
77) Benzidine	19.445	184	526463	52.499	ng	100
78) Pyrene	19.621	202	5467448	64.482	ng	99
80) Butylbenzylphthalate	20.539	149	1822716	69.083	ng	91
81) Benzo(a)anthracene	21.433	228	5054698	65.863	ng	100
82) 3,3'-Dichlorobenzidine	21.350	252	1313089	63.583	ng	100
83) Chrysene	21.492	228	4675833	65.030	ng	100
84) Bis(2-ethylhexyl)phtha...	21.380	149	2998277	66.289	ng	100
85) Di-n-octyl phthalate	22.544	149	4772040	66.876	ng	99
87) Indeno(1,2,3-cd)pyrene	27.909	276	3473188	68.678	ng	99
88) Benzo(b)fluoranthene	23.527	252	4312843	67.700	ng	100
89) Benzo(k)fluoranthene	23.591	252	4204462	67.952	ng	99
90) Benzo(a)pyrene	24.338	252	3412806	67.601	ng	99
91) Dibenzo(a,h)anthracene	27.974	278	2781271	69.585	ng	98
92) Benzo(g,h,i)perylene	28.973	276	3012226	67.465	ng	99

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

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