

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM070925\
 Data File : BM050384.D
 Acq On : 08 Jul 2025 17:22
 Operator : RC/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC080

Quant Time: Jul 08 18:24:43 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM070925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 08 17:58:12 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.869	152	429681	20.000	ng	0.00
21) Naphthalene-d8	10.663	136	1643859	20.000	ng	0.00
39) Acenaphthene-d10	14.498	164	1043719	20.000	ng	0.00
64) Phenanthrene-d10	17.227	188	1944455	20.000	ng	0.00
76) Chrysene-d12	21.457	240	2055292	20.000	ng	0.00
86) Perylene-d12	24.498	264	1953481	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.440	112	4014825	160.835	ng	0.00
7) Phenol-d6	7.028	99	5149049	163.630	ng	0.00
23) Nitrobenzene-d5	9.028	82	5264647	163.391	ng	0.00
42) 2,4,6-Tribromophenol	15.980	330	2335664	184.182	ng	0.00
45) 2-Fluorobiphenyl	13.122	172	13148691	155.576	ng	0.00
79) Terphenyl-d14	19.839	244	14435460	123.577	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.334	88	785309	76.507	ng	99
3) Pyridine	3.734	79	2123609	78.804	ng	98
4) n-Nitrosodimethylamine	3.640	42	970223	77.228	ng	96
6) Aniline	7.193	93	3301305	81.748	ng	99
8) 2-Chlorophenol	7.428	128	2162654	82.163	ng	99
10) Phenol	7.052	94	2637766	80.368	ng	100
11) bis(2-Chloroethyl)ether	7.287	93	2092911	79.625	ng	99
12) 1,3-Dichlorobenzene	7.757	146	2517372	78.654	ng	98
13) 1,4-Dichlorobenzene	7.904	146	2555213	78.187	ng	99
14) 1,2-Dichlorobenzene	8.222	146	2454039	78.434	ng	99
15) Benzyl Alcohol	8.104	79	1807870	81.539	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.399	45	2991629	77.270	ng	99
17) 2-Methylphenol	8.304	107	1732716	81.397	ng	99
18) Hexachloroethane	8.951	117	932539	81.135	ng	99
19) n-Nitroso-di-n-propyla...	8.681	70	1587587	83.281	ng	95
20) 3+4-Methylphenols	8.634	107	2300155	80.491	ng	100
22) Acetophenone	8.687	105	3217324	78.279	ng	98
24) Nitrobenzene	9.063	77	2296616	79.875	ng	100
25) Isophorone	9.598	82	4289620	82.044	ng	100
26) 2-Nitrophenol	9.775	139	1095059	93.448	ng	97
27) 2,4-Dimethylphenol	9.834	122	2026460	81.267	ng	97
28) bis(2-Chloroethoxy)met...	10.075	93	2760921	79.691	ng	99
29) 2,4-Dichlorophenol	10.310	162	2134843	83.334	ng	98
30) 1,2,4-Trichlorobenzene	10.528	180	2483928	81.052	ng	99
31) Naphthalene	10.716	128	6520664	78.095	ng	100
32) Benzoic acid	9.993	122	1311597	81.001	ng	97
33) 4-Chloroaniline	10.816	127	2873151	81.395	ng	99
34) Hexachlorobutadiene	11.010	225	1546563	82.684	ng	100
35) Caprolactam	11.604	113	594080	84.949	ng	93
36) 4-Chloro-3-methylphenol	11.940	107	1983864	82.929	ng	100
37) 2-Methylnaphthalene	12.322	142	4254768	80.705	ng	99
38) 1-Methylnaphthalene	12.545	142	4475032	80.207	ng	99
40) 1,2,4,5-Tetrachloroben...	12.692	216	2824950	85.112	ng	99
41) Hexachlorocyclopentadiene	12.681	237	1914528	90.239	ng	99
43) 2,4,6-Trichlorophenol	12.928	196	1782842	87.263	ng	99
44) 2,4,5-Trichlorophenol	12.998	196	1919360	86.021	ng	97

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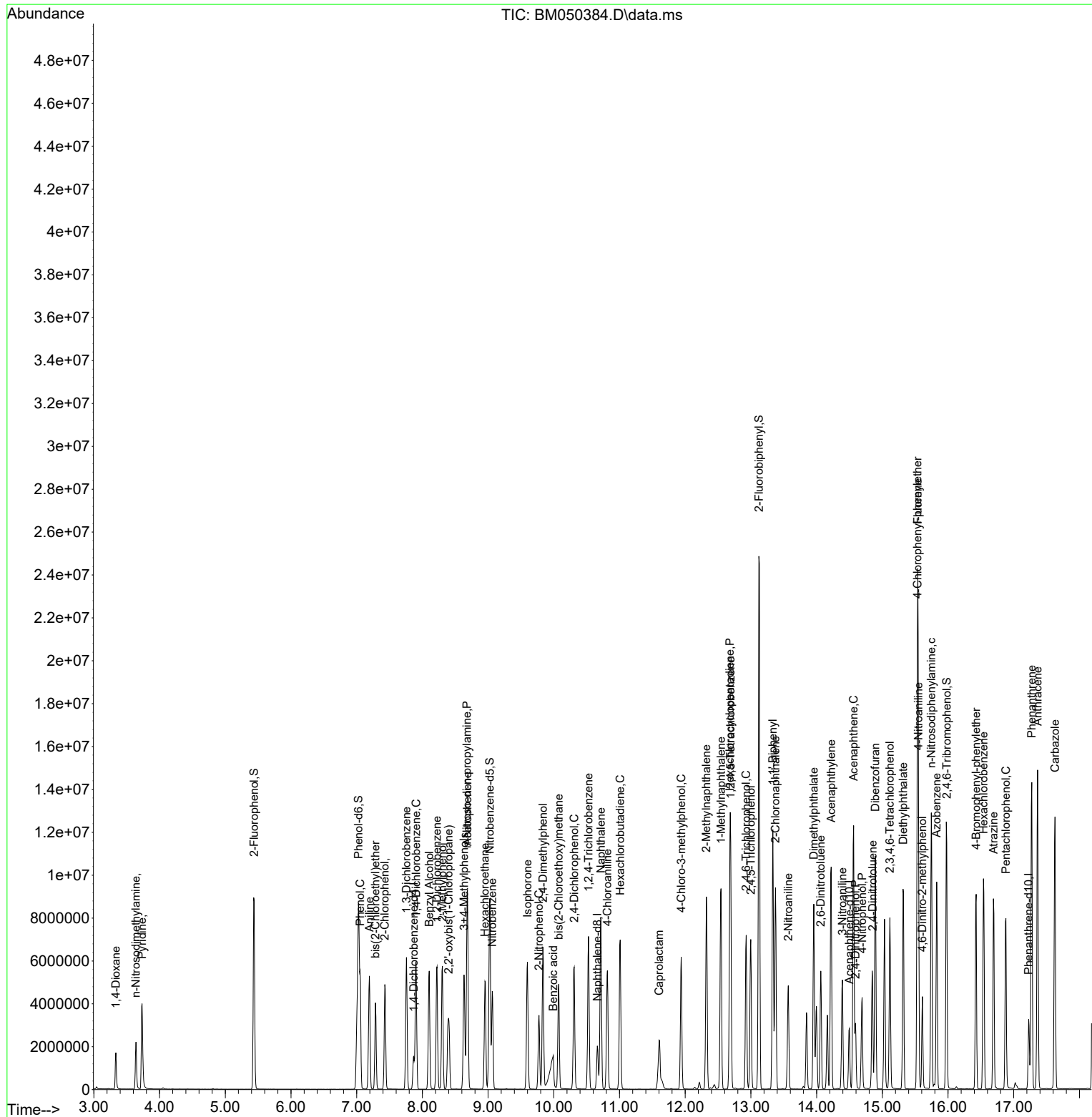
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 1,1'-Biphenyl	13.334	154	6141549	79.310	ng	100
47) 2-Chloronaphthalene	13.375	162	4889665	79.196	ng	99
48) 2-Nitroaniline	13.569	65	1233222	89.485	ng	96
49) Acenaphthylene	14.222	152	7559726	81.019	ng	100
50) Dimethylphthalate	13.957	163	5732691	79.779	ng	100
51) 2,6-Dinitrotoluene	14.063	165	1207179	87.496	ng	99
52) Acenaphthene	14.563	154	4833886	81.486	ng	99
53) 3-Nitroaniline	14.392	138	1297724	88.444	ng	99
54) 2,4-Dinitrophenol	14.592	184	626569	82.116	ng	86
55) Dibenzofuran	14.892	168	7143343	78.152	ng	99
56) 4-Nitrophenol	14.692	139	1075848	85.262	ng	93
57) 2,4-Dinitrotoluene	14.851	165	1674329	79.189	ng	94
58) Fluorene	15.539	166	6138107	81.530	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.116	232	1636646	87.136	ng	97
60) Diethylphthalate	15.316	149	5511226	80.291	ng	100
61) 4-Chlorophenyl-phenyle...	15.533	204	3298863	83.798	ng	98
62) 4-Nitroaniline	15.551	138	1336300	77.536	ng	97
63) Azobenzene	15.828	77	5026045	79.741	ng	99
65) 4,6-Dinitro-2-methylph...	15.610	198	912781	82.140	ng	92
66) n-Nitrosodiphenylamine	15.745	169	5086187	83.592	ng	100
67) 4-Bromophenyl-phenylether	16.427	248	1878071	87.664	ng	97
68) Hexachlorobenzene	16.539	284	2164238	85.534	ng	97
69) Atrazine	16.692	200	1767182	88.450	ng	99
70) Pentachlorophenol	16.880	266	1425059	90.473	ng	99
71) Phenanthrene	17.274	178	8893346	80.828	ng	99
72) Anthracene	17.363	178	9167986	83.709	ng	99
73) Carbazole	17.627	167	8180242	82.547	ng	99
74) Di-n-butylphthalate	18.198	149	9300935	85.749	ng	99
75) Fluoranthene	19.280	202	10317307	86.256	ng	99
77) Benzidine	19.457	184	3817410	72.819	ng	100
78) Pyrene	19.639	202	10818153	82.185	ng	99
80) Butylbenzylphthalate	20.539	149	3938114	90.734	ng	97
81) Benzo(a)anthracene	21.439	228	11055858	82.558	ng	99
82) 3,3'-Dichlorobenzidine	21.362	252	3825108	84.699	ng	99
83) Chrysene	21.504	228	10270454	81.310	ng	99
84) Bis(2-ethylhexyl)phtha...	21.374	149	6100517	88.540	ng	# 96
85) Di-n-octyl phthalate	22.527	149	9722259	90.093	ng	99
87) Indeno(1,2,3-cd)pyrene	27.956	276	12120915	87.270	ng	99
88) Benzo(b)fluoranthene	23.545	252	10371642	86.316	ng	99
89) Benzo(k)fluoranthene	23.609	252	10546917	84.592	ng	100
90) Benzo(a)pyrene	24.362	252	9879526	87.833	ng	100
91) Dibenzo(a,h)anthracene	28.015	278	10165335	86.905	ng	99
92) Benzo(g,h,i)perylene	29.027	276	9504954	85.979	ng	99

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

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