

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041525\
 Data File : BM049929.D
 Acq On : 15 Apr 2025 11:15
 Operator : RC/JU
 Sample : PB167577BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB167577BS

Quant Time: Apr 15 12:36:55 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM040825.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 09 04:00:55 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	372144	20.000	ng	-0.01
21) Naphthalene-d8	10.569	136	1305000	20.000	ng	0.00
39) Acenaphthene-d10	14.416	164	859595	20.000	ng	0.00
64) Phenanthrene-d10	17.162	188	1699421	20.000	ng	0.00
76) Chrysene-d12	21.403	240	1667631	20.000	ng	0.00
86) Perylene-d12	24.403	264	1728437	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.363	112	2800461	127.784	ng	0.00
7) Phenol-d6	6.951	99	3447148	126.422	ng	0.00
23) Nitrobenzene-d5	8.928	82	1976870	84.354	ng	0.00
42) 2,4,6-Tribromophenol	15.910	330	1754313	140.310	ng	0.00
45) 2-Fluorobiphenyl	13.039	172	5335402	84.166	ng	0.00
79) Terphenyl-d14	19.786	244	8761319	98.458	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.257	88	329364	36.492	ng	97
3) Pyridine	3.657	79	902568	38.061	ng	100
4) n-Nitrosodimethylamine	3.563	42	383854	42.535	ng	99
6) Aniline	7.098	93	621845	26.615	ng	99
8) 2-Chlorophenol	7.340	128	1033856	46.032	ng	99
9) Benzaldehyde	6.910	77	329066	23.517	ng	98
10) Phenol	6.981	94	1231858	46.295	ng	99
11) bis(2-Chloroethyl)ether	7.192	93	988233	47.360	ng	100
12) 1,3-Dichlorobenzene	7.663	146	1168508	44.006	ng	99
13) 1,4-Dichlorobenzene	7.804	146	1172347	43.879	ng	100
14) 1,2-Dichlorobenzene	8.122	146	1132578	45.005	ng	99
15) Benzyl Alcohol	8.016	79	764123	44.503	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.298	45	1103982	47.468	ng	99
17) 2-Methylphenol	8.228	107	800269	48.804	ng	99
18) Hexachloroethane	8.851	117	410036	44.112	ng	94
19) n-Nitroso-di-n-propyla...	8.581	70	690538	45.725	ng	99
20) 3+4-Methylphenols	8.551	107	1068101	47.778	ng	100
22) Acetophenone	8.592	105	1379423	43.494	ng	100
24) Nitrobenzene	8.969	77	989812	43.863	ng	98
25) Isophorone	9.498	82	1880139	47.891	ng	99
26) 2-Nitrophenol	9.681	139	536665	44.870	ng	99
27) 2,4-Dimethylphenol	9.751	122	893806	65.942	ng	99
28) bis(2-Chloroethoxy)met...	9.975	93	1187922	45.197	ng	99
29) 2,4-Dichlorophenol	10.228	162	1028800	45.661	ng	99
30) 1,2,4-Trichlorobenzene	10.433	180	1143710	43.879	ng	99
31) Naphthalene	10.616	128	2836214	42.296	ng	99
32) Benzoic acid	9.904	122	593692m	37.589	ng	
33) 4-Chloroaniline	10.739	127	573155	24.206	ng	99
34) Hexachlorobutadiene	10.904	225	720159	44.667	ng	98
35) Caprolactam	11.516	113	284818	44.075	ng	96
36) 4-Chloro-3-methylphenol	11.869	107	908816	46.048	ng	100
37) 2-Methylnaphthalene	12.233	142	1891326	40.032	ng	100
38) 1-Methylnaphthalene	12.451	142	1994230	43.326	ng	99
40) 1,2,4,5-Tetrachloroben...	12.604	216	1310150	43.566	ng	99
41) Hexachlorocyclopentadiene	12.586	237	1722388	163.451	ng	100
43) 2,4,6-Trichlorophenol	12.851	196	877590	45.860	ng	98

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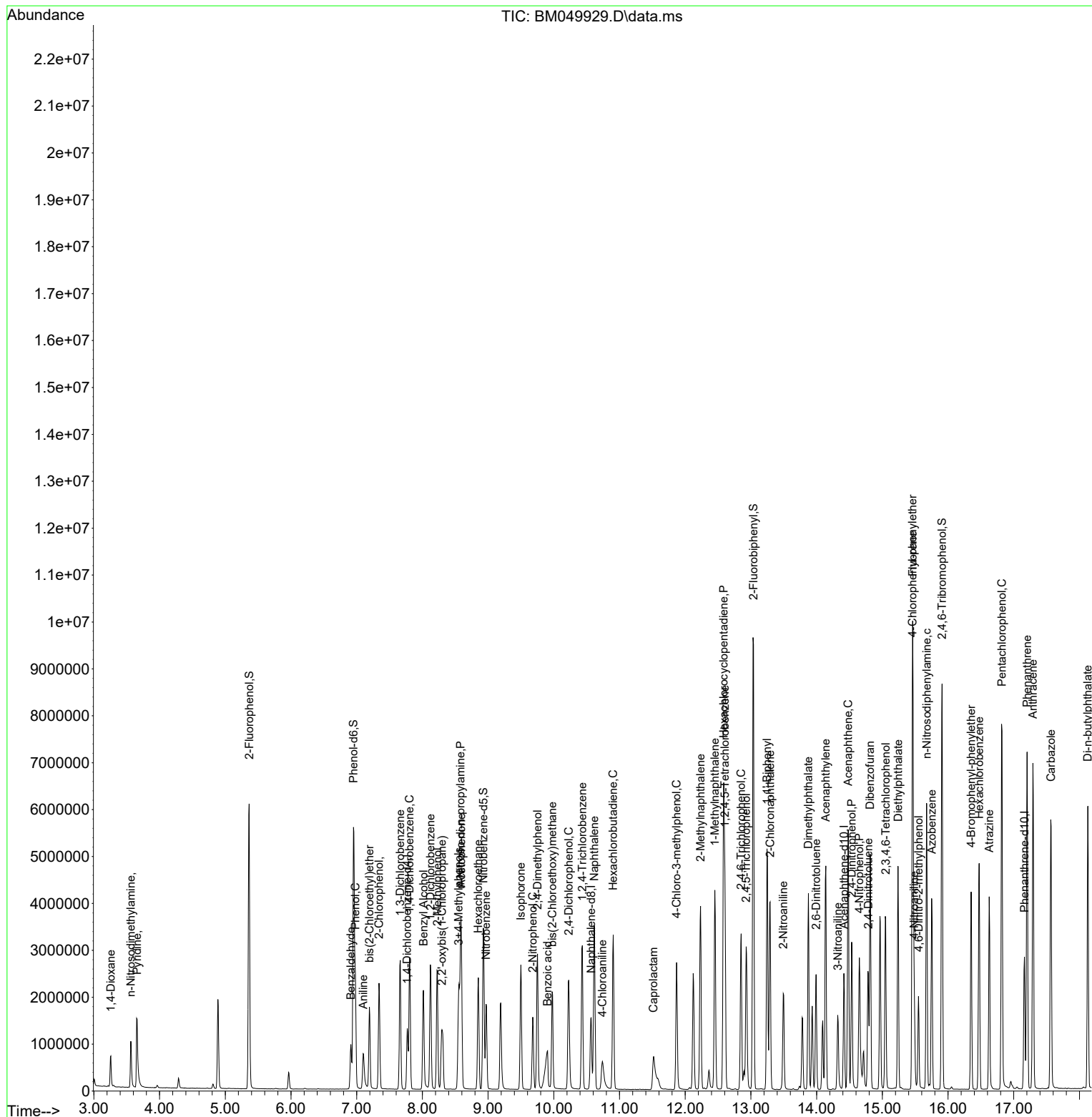
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.927	196	951142	45.738	ng	99
46) 1,1'-Biphenyl	13.245	154	2737950	42.848	ng	100
47) 2-Chloronaphthalene	13.292	162	2184876	43.503	ng	99
48) 2-Nitroaniline	13.492	65	545085	46.913	ng	96
49) Acenaphthylene	14.139	152	3457932	47.266	ng	100
50) Dimethylphthalate	13.874	163	2718514	44.828	ng	99
51) 2,6-Dinitrotoluene	13.992	165	589963	46.433	ng	97
52) Acenaphthene	14.480	154	2002580	43.034	ng	99
53) 3-Nitroaniline	14.321	138	466945	38.025	ng	96
54) 2,4-Dinitrophenol	14.533	184	796405	85.963	ng	97
55) Dibenzofuran	14.816	168	3286910	42.252	ng	99
56) 4-Nitrophenol	14.651	139	1021173	93.317	ng	98
57) 2,4-Dinitrotoluene	14.786	165	834436	49.229	ng	95
58) Fluorene	15.463	166	2755356	44.559	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.051	232	819934	44.985	ng	98
60) Diethylphthalate	15.239	149	2539531	43.668	ng	98
61) 4-Chlorophenyl-phenyle...	15.457	204	1515550	45.714	ng	99
62) 4-Nitroaniline	15.486	138	573366	45.150	ng	100
63) Azobenzene	15.751	77	2135715	42.894	ng	98
65) 4,6-Dinitro-2-methylph...	15.551	198	497708	46.476	ng	98
66) n-Nitrosodiphenylamine	15.674	169	2348681	46.182	ng	100
67) 4-Bromophenyl-phenylether	16.351	248	913125	46.285	ng	97
68) Hexachlorobenzene	16.474	284	1053168	46.541	ng	99
69) Atrazine	16.627	200	816421	61.914	ng	98
70) Pentachlorophenol	16.821	266	1497814	89.906	ng	99
71) Phenanthrene	17.204	178	4249848	45.614	ng	100
72) Anthracene	17.292	178	4334490	47.208	ng	100
73) Carbazole	17.568	167	3895183	45.042	ng	99
74) Di-n-butylphthalate	18.127	149	4400587	44.179	ng	100
75) Fluoranthene	19.221	202	5134794	44.824	ng	99
77) Benzidine	19.409	184	1800665	81.392	ng	100
78) Pyrene	19.586	202	5314431	46.241	ng	99
80) Butylbenzylphthalate	20.480	149	1979147	45.829	ng	99
81) Benzo(a)anthracene	21.380	228	5216434	47.067	ng	100
82) 3,3'-Dichlorobenzidine	21.303	252	1250641	29.881	ng	100
83) Chrysene	21.445	228	4810031	45.650	ng	100
84) Bis(2-ethylhexyl)phtha...	21.303	149	2821380	44.841	ng	98
85) Di-n-octyl phthalate	22.439	149	4860081	44.153	ng	100
87) Indeno(1,2,3-cd)pyrene	27.803	276	5958306	46.246	ng	99
88) Benzo(b)fluoranthene	23.456	252	4896246	44.355	ng	99
89) Benzo(k)fluoranthene	23.521	252	4979762	46.551	ng	100
90) Benzo(a)pyrene	24.262	252	4746443	48.931	ng	99
91) Dibenzo(a,h)anthracene	27.850	278	4886859	46.049	ng	99
92) Benzo(g,h,i)perylene	28.856	276	4718483	43.511	ng	98

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

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