

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020525\
 Data File : BM049565.D
 Acq On : 05 Feb 2025 11:44
 Operator : RC/JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC005

Quant Time: Feb 06 04:37:42 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	281355	20.000	ng	0.00
21) Naphthalene-d8	10.622	136	1040421	20.000	ng	0.00
39) Acenaphthene-d10	14.463	164	663031	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	1282078	20.000	ng	0.00
76) Chrysene-d12	21.445	240	1050642	20.000	ng	0.00
86) Perylene-d12	24.474	264	939392	20.000	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	164670	9.412	ng	0.00
7) Phenol-d6	6.975	99	207416	9.049	ng	0.00
23) Nitrobenzene-d5	8.975	82	177204	8.761	ng	0.00
42) 2,4,6-Tribromophenol	0.000	330	0d	0.000	ng	
45) 2-Fluorobiphenyl	13.092	172	433007	8.460	ng	0.00
79) Terphenyl-d14	19.833	244	552192	8.026	ng	0.00

Target Compounds					Qvalue
2) 1,4-Dioxane	3.293	88	35564	4.834	ng 95
3) Pyridine	3.693	79	98511	4.819	ng 96
4) n-Nitrosodimethylamine	3.599	42	44337	4.636	ng # 96
6) Aniline	7.146	93	102752	4.730	ng 98
8) 2-Chlorophenol	7.387	128	79936	4.576	ng 96
9) Benzaldehyde	6.957	77	67388	5.552	ng 99
10) Phenol	7.004	94	104865	4.645	ng 97
11) bis(2-Chloroethyl)ether	7.246	93	86785	4.751	ng 96
12) 1,3-Dichlorobenzene	7.716	146	98691	4.867	ng 96
13) 1,4-Dichlorobenzene	7.863	146	99490	4.850	ng 99
14) 1,2-Dichlorobenzene	8.175	146	95387	4.816	ng 98
15) Benzyl Alcohol	8.057	79	72503	4.546	ng 100
16) 2,2'-oxybis(1-Chloropr...	8.363	45	117291	4.770	ng 97
17) 2-Methylphenol	8.257	107	64889	4.700	ng 98
18) Hexachloroethane	8.910	117	35779	4.775	ng 96
19) n-Nitroso-di-n-propyla...	8.628	70	69534	4.632	ng 98
20) 3+4-Methylphenols	8.581	107	82241	4.513	ng 97
22) Acetophenone	8.640	105	130422	4.654	ng 98
24) Nitrobenzene	9.016	77	88390	4.498	ng 98
25) Isophorone	9.545	82	167818	4.600	ng 100
26) 2-Nitrophenol	9.734	139	20705	5.598	ng 97
27) 2,4-Dimethylphenol	9.792	122	62122	5.386	ng 96
28) bis(2-Chloroethoxy)met...	10.034	93	110852	4.718	ng 99
29) 2,4-Dichlorophenol	10.263	162	66638	4.195	ng 98
30) 1,2,4-Trichlorobenzene	10.486	180	91662	4.710	ng 98
31) Naphthalene	10.675	128	256675	4.678	ng 99
33) 4-Chloroaniline	10.769	127	96388	4.693	ng 99
34) Hexachlorobutadiene	10.975	225	53346	4.577	ng 98
35) Caprolactam	11.504	113	19856	4.097	ng 94
36) 4-Chloro-3-methylphenol	11.892	107	67154	4.239	ng 98
37) 2-Methylnaphthalene	12.286	142	176822	4.577	ng 99
38) 1-Methylnaphthalene	12.504	142	173149	4.606	ng 100
40) 1,2,4,5-Tetrachloroben...	12.657	216	95516	4.382	ng 99
41) Hexachlorocyclopentadiene	12.645	237	20613	3.704	ng 97
43) 2,4,6-Trichlorophenol	12.892	196	45994	3.866	ng 98
44) 2,4,5-Trichlorophenol	12.957	196	47890	3.663	ng 96

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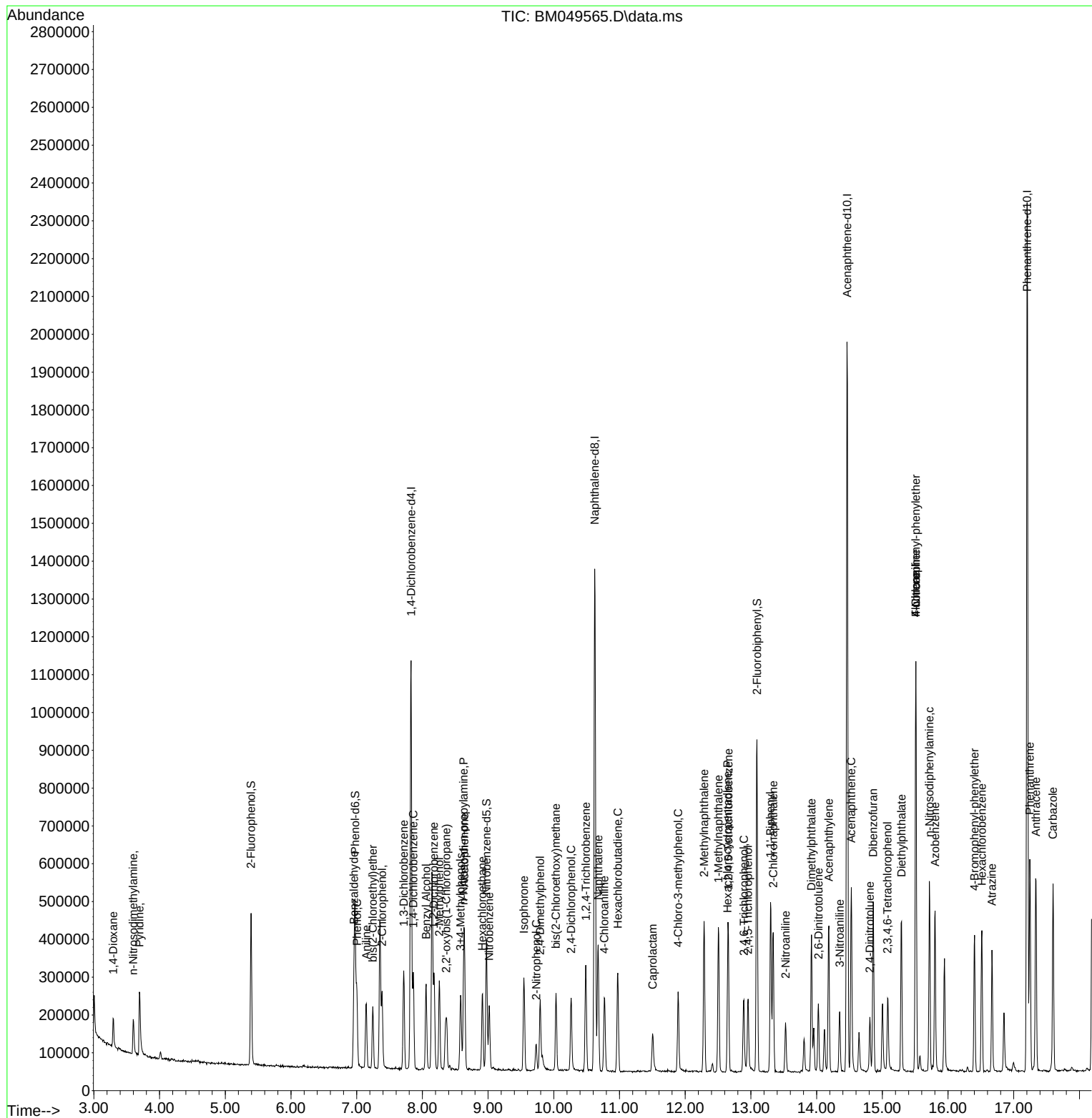
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 1,1'-Biphenyl	13.298	154	229625	4.540	ng	98
47) 2-Chloronaphthalene	13.339	162	177515	4.536	ng	98
48) 2-Nitroaniline	13.527	65	31147	6.886	ng	91
49) Acenaphthylene	14.186	152	262019	4.541	ng	99
50) Dimethylphthalate	13.922	163	220396	4.629	ng	99
51) 2,6-Dinitrotoluene	14.027	165	28514	3.461	ng	85
52) Acenaphthene	14.527	154	166637	4.351	ng	98
53) 3-Nitroaniline	14.351	138	29903	3.565	ng #	91
55) Dibenzofuran	14.863	168	277206	4.545	ng	100
57) 2,4-Dinitrotoluene	14.810	165	29729	5.525	ng	88
58) Fluorene	15.510	166	216619	4.152	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.080	232	40263	3.646	ng	96
60) Diethylphthalate	15.292	149	215092	4.495	ng	99
61) 4-Chlorophenyl-phenyle...	15.510	204	110799	3.990	ng	92
62) 4-Nitroaniline	15.510	138	31375	6.847	ng #	82
63) Azobenzene	15.804	77	237223	4.749	ng	98
66) n-Nitrosodiphenylamine	15.716	169	176950	4.376	ng	100
67) 4-Bromophenyl-phenylether	16.404	248	63592	4.269	ng	96
68) Hexachlorobenzene	16.516	284	74533	4.360	ng	96
69) Atrazine	16.668	200	52467	4.365	ng	99
71) Phenanthrene	17.245	178	318001	4.354	ng	99
72) Anthracene	17.339	178	311738	4.311	ng	99
73) Carbazole	17.598	167	289997	4.569	ng	98
74) Di-n-butylphthalate	18.186	149	333555	4.141	ng	99
75) Fluoranthene	19.262	202	346582	4.053	ng	99
77) Benzidine	19.439	184	54625	5.871	ng	99
78) Pyrene	19.621	202	354809	4.510	ng	98
80) Butylbenzylphthalate	20.539	149	87083	3.557	ng	99
81) Benzo(a)anthracene	21.427	228	312451	4.388	ng	99
82) 3,3'-Dichlorobenzidine	21.350	252	80259	4.189	ng	99
83) Chrysene	21.486	228	291003	4.362	ng	99
84) Bis(2-ethylhexyl)phtha...	21.386	149	171121	4.078	ng	99
85) Di-n-octyl phthalate	22.545	149	265303	4.007	ng	97
87) Indeno(1,2,3-cd)pyrene	27.891	276	204636	3.989	ng	99
88) Benzo(b)fluoranthene	23.521	252	267263	4.136	ng	98
89) Benzo(k)fluoranthene	23.580	252	258722	4.122	ng	99
90) Benzo(a)pyrene	24.333	252	210395	4.108	ng	98
91) Dibenzo(a,h)anthracene	27.968	278	157949	3.896	ng	98
92) Benzo(g,h,i)perylene	28.950	276	186220	4.111	ng	98

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

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