

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020625\
 Data File : BM049594.D
 Acq On : 07 Feb 2025 00:52
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Feb 07 01:34:52 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 04:55:28 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	252484	20.000	ng	0.00
21) Naphthalene-d8	10.622	136	937146	20.000	ng	0.00
39) Acenaphthene-d10	14.463	164	583935	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	1104130	20.000	ng	0.00
76) Chrysene-d12	21.445	240	1124649	20.000	ng	0.00
86) Perylene-d12	24.474	264	936920	20.000	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	1285109	81.849	ng	0.00
7) Phenol-d6	6.981	99	1677721	81.561	ng	0.00
23) Nitrobenzene-d5	8.981	82	1577845	86.603	ng	0.00
42) 2,4,6-Tribromophenol	15.945	330	574481	83.028	ng	0.00
45) 2-Fluorobiphenyl	13.092	172	3887755	86.249	ng	0.00
79) Terphenyl-d14	19.833	244	6699385	90.964	ng	0.00

Target Compounds					Qvalue
2) 1,4-Dioxane	3.299	88	274955	41.650	ng 100
3) Pyridine	3.693	79	752603	41.025	ng 100
4) n-Nitrosodimethylamine	3.599	42	356200	41.501	ng 97
6) Aniline	7.145	93	759049	38.934	ng 98
8) 2-Chlorophenol	7.387	128	626642	39.970	ng 98
9) Benzaldehyde	6.957	77	468038	42.972	ng 99
10) Phenol	7.010	94	817552	40.352	ng 98
11) bis(2-Chloroethyl)ether	7.245	93	663825	40.499	ng 96
12) 1,3-Dichlorobenzene	7.716	146	743968	40.882	ng 98
13) 1,4-Dichlorobenzene	7.857	146	754363	40.983	ng 99
14) 1,2-Dichlorobenzene	8.175	146	731707	41.169	ng 100
15) Benzyl Alcohol	8.057	79	557993	38.991	ng 98
16) 2,2'-oxybis(1-Chloropr...	8.363	45	914224	41.433	ng 97
17) 2-Methylphenol	8.257	107	501690	40.497	ng 97
18) Hexachloroethane	8.910	117	278324	41.396	ng 97
19) n-Nitroso-di-n-propyla...	8.634	70	555103	41.210	ng 99
20) 3+4-Methylphenols	8.587	107	668477	40.876	ng 98
22) Acetophenone	8.639	105	1042353	41.296	ng # 97
24) Nitrobenzene	9.022	77	760656	42.972	ng 100
25) Isophorone	9.545	82	1306320	39.757	ng 100
26) 2-Nitrophenol	9.734	139	214593	37.484	ng 98
27) 2,4-Dimethylphenol	9.792	122	394781	37.999	ng 97
28) bis(2-Chloroethoxy)met...	10.034	93	861535	40.711	ng 99
29) 2,4-Dichlorophenol	10.263	162	590841	41.295	ng 98
30) 1,2,4-Trichlorobenzene	10.486	180	709167	40.456	ng 97
31) Naphthalene	10.675	128	2002909	40.524	ng 99
32) Benzoic acid	9.910	122	225319	38.695	ng 95
33) 4-Chloroaniline	10.769	127	722271	39.042	ng 99
34) Hexachlorobutadiene	10.969	225	434622	41.403	ng 100
35) Caprolactam	11.539	113	171049	39.187	ng 99
36) 4-Chloro-3-methylphenol	11.898	107	575286	40.318	ng 97
37) 2-Methylnaphthalene	12.286	142	1393084	40.032	ng 98
38) 1-Methylnaphthalene	12.504	142	1339011	39.548	ng 100
40) 1,2,4,5-Tetrachloroben...	12.651	216	801288	41.736	ng 99
41) Hexachlorocyclopentadiene	12.645	237	191115	38.994	ng 99
43) 2,4,6-Trichlorophenol	12.892	196	419910	40.081	ng 97

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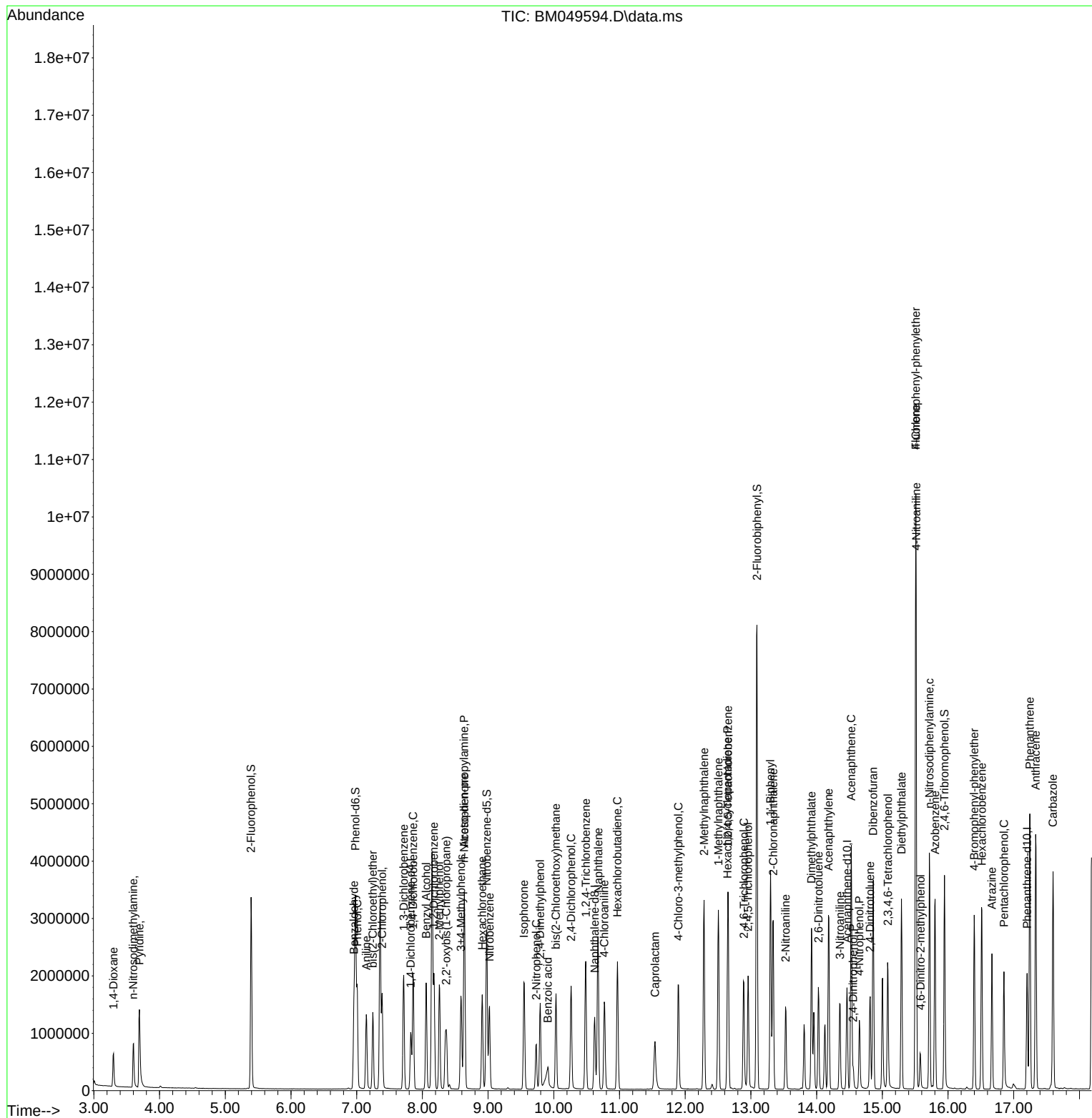
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.957	196	479844	41.678	ng	96
46) 1,1'-Biphenyl	13.298	154	1804680	40.513	ng	99
47) 2-Chloronaphthalene	13.339	162	1390026	40.331	ng	99
48) 2-Nitroaniline	13.527	65	356049	39.940	ng	94
49) Acenaphthylene	14.180	152	2044800	40.235	ng	100
50) Dimethylphthalate	13.921	163	1667137	39.755	ng	99
51) 2,6-Dinitrotoluene	14.027	165	320528	44.172	ng	93
52) Acenaphthene	14.527	154	1379788	40.903	ng	99
53) 3-Nitroaniline	14.357	138	323200	43.747	ng	99
54) 2,4-Dinitrophenol	14.557	184	62516	38.254	ng	# 87
55) Dibenzofuran	14.863	168	2183999	40.661	ng	100
56) 4-Nitrophenol	14.651	139	252332	40.517	ng	99
57) 2,4-Dinitrotoluene	14.816	165	416433	42.335	ng	99
58) Fluorene	15.510	166	2026194	44.096	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.086	232	410589	42.211	ng	100
60) Diethylphthalate	15.292	149	1697401	40.276	ng	100
61) 4-Chlorophenyl-phenyle...	15.510	204	1085560	44.387	ng	99
62) 4-Nitroaniline	15.515	138	392660	41.679	ng	98
63) Azobenzene	15.804	77	1826054	41.511	ng	99
65) 4,6-Dinitro-2-methylph...	15.580	198	106790	37.603	ng	99
66) n-Nitrosodiphenylamine	15.715	169	1415298	40.645	ng	99
67) 4-Bromophenyl-phenylether	16.398	248	516640	40.277	ng	98
68) Hexachlorobenzene	16.515	284	599954	40.751	ng	96
69) Atrazine	16.668	200	374530	36.183	ng	97
70) Pentachlorophenol	16.851	266	312275	36.109	ng	99
71) Phenanthrene	17.245	178	2584326	41.085	ng	99
72) Anthracene	17.333	178	2567637	41.226	ng	100
73) Carbazole	17.598	167	2223107	40.668	ng	100
74) Di-n-butylphthalate	18.186	149	2792159	40.254	ng	100
75) Fluoranthene	19.256	202	3100217	42.098	ng	100
77) Benzidine	19.439	184	300777	30.200	ng	98
78) Pyrene	19.621	202	3208842	38.105	ng	99
80) Butylbenzylphthalate	20.533	149	888399	33.904	ng	96
81) Benzo(a)anthracene	21.427	228	3128202	41.042	ng	100
82) 3,3'-Dichlorobenzidine	21.345	252	792327	38.631	ng	99
83) Chrysene	21.486	228	2966738	41.545	ng	100
84) Bis(2-ethylhexyl)phtha...	21.374	149	1660616	36.968	ng	99
85) Di-n-octyl phthalate	22.533	149	2428980	34.275	ng	99
87) Indeno(1,2,3-cd)pyrene	27.897	276	2197229	42.942	ng	100
88) Benzo(b)fluoranthene	23.521	252	2660341	41.275	ng	99
89) Benzo(k)fluoranthene	23.586	252	2659172	42.477	ng	99
90) Benzo(a)pyrene	24.333	252	2126485	41.632	ng	99
91) Dibenzo(a,h)anthracene	27.962	278	1742460	43.088	ng	97
92) Benzo(g,h,i)perylene	28.962	276	1938625	42.914	ng	99

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

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