

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040825\
 Data File : BM049851.D
 Acq On : 08 Apr 2025 15:32
 Operator : RC/JU
 Sample : SSTDICC020
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC020

Quant Time: Apr 09 03:01:01 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 02:53:19 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	318059	20.000	ng	0.00
21) Naphthalene-d8	10.575	136	1116195	20.000	ng	0.00
39) Acenaphthene-d10	14.422	164	702854	20.000	ng	0.00
64) Phenanthrene-d10	17.163	188	1378813	20.000	ng	0.00
76) Chrysene-d12	21.403	240	1339758	20.000	ng	0.00
86) Perylene-d12	24.403	264	1486954	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.363	112	771430	41.186	ng	0.00
7) Phenol-d6	6.951	99	968531	41.560	ng	0.00
23) Nitrobenzene-d5	8.934	82	817733	40.795	ng	0.00
42) 2,4,6-Tribromophenol	15.910	330	393158	38.457	ng	0.00
45) 2-Fluorobiphenyl	13.039	172	2107153	40.653	ng	0.00
79) Terphenyl-d14	19.792	244	3051000	42.677	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.263	88	159450	20.670	ng	99
3) Pyridine	3.663	79	420529	20.757	ng	99
4) n-Nitrosodimethylamine	3.569	42	157731	20.450	ng	96
6) Aniline	7.104	93	443873	22.228	ng	100
8) 2-Chlorophenol	7.346	128	397524	20.710	ng	99
9) Benzaldehyde	6.916	77	266386	22.294	ng	99
10) Phenol	6.975	94	471508	20.733	ng	99
11) bis(2-Chloroethyl)ether	7.198	93	364893	20.461	ng	98
12) 1,3-Dichlorobenzene	7.669	146	469124	20.672	ng	100
13) 1,4-Dichlorobenzene	7.816	146	473136	20.720	ng	100
14) 1,2-Dichlorobenzene	8.134	146	446077	20.740	ng	99
15) Benzyl Alcohol	8.016	79	302185	20.592	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.310	45	417355	20.996	ng	99
17) 2-Methylphenol	8.222	107	291405	20.793	ng	99
18) Hexachloroethane	8.857	117	165304	20.807	ng	98
19) n-Nitroso-di-n-propyla...	8.581	70	273519	21.191	ng	99
20) 3+4-Methylphenols	8.551	107	396422	20.748	ng	98
22) Acetophenone	8.598	105	554070	20.425	ng	# 99
24) Nitrobenzene	8.975	77	395594	20.496	ng	99
25) Isophorone	9.504	82	683707	20.361	ng	100
26) 2-Nitrophenol	9.687	139	203797	19.922	ng	97
27) 2,4-Dimethylphenol	9.751	122	231983	20.010	ng	100
28) bis(2-Chloroethoxy)met...	9.981	93	459217	20.427	ng	99
29) 2,4-Dichlorophenol	10.222	162	388584	20.164	ng	99
30) 1,2,4-Trichlorobenzene	10.439	180	448099	20.099	ng	100
31) Naphthalene	10.622	128	1166673	20.341	ng	99
32) Benzoic acid	9.839	122	195350	19.812	ng	97
33) 4-Chloroaniline	10.728	127	420078	20.742	ng	99
34) Hexachlorobutadiene	10.916	225	273841	19.857	ng	99
35) Caprolactam	11.504	113	111645	20.199	ng	98
36) 4-Chloro-3-methylphenol	11.863	107	340457	20.168	ng	98
37) 2-Methylnaphthalene	12.239	142	813541	20.132	ng	100
38) 1-Methylnaphthalene	12.457	142	793590	20.158	ng	99
40) 1,2,4,5-Tetrachloroben...	12.610	216	489705	19.915	ng	99
41) Hexachlorocyclopentadiene	12.592	237	173683	20.158	ng	99
43) 2,4,6-Trichlorophenol	12.851	196	311742	19.924	ng	97

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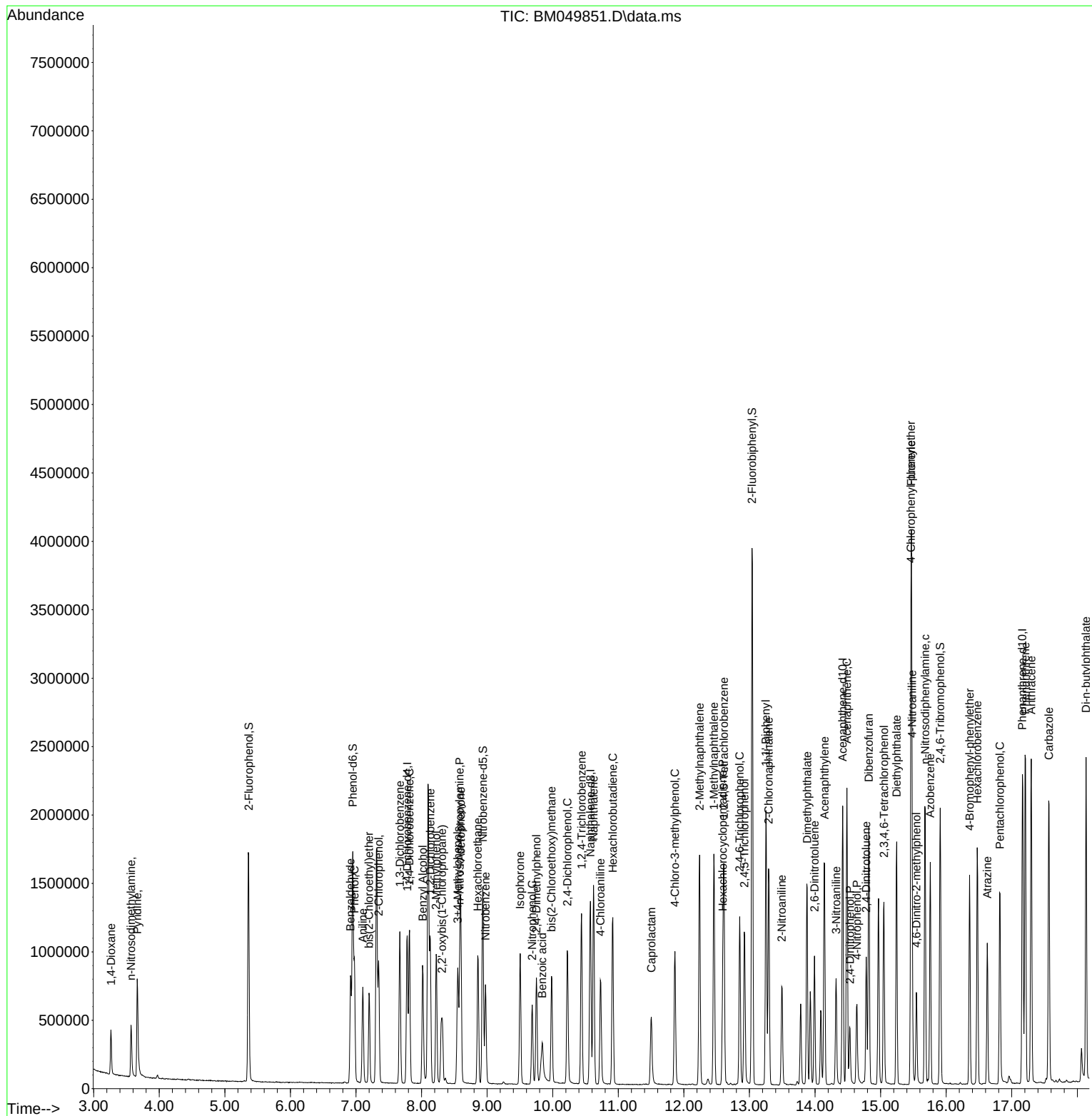
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.922	196	332054	19.528	ng	96
46) 1,1'-Biphenyl	13.251	154	1063836	20.361	ng	99
47) 2-Chloronaphthalene	13.292	162	835052	20.334	ng	100
48) 2-Nitroaniline	13.492	65	191426	20.149	ng	94
49) Acenaphthylene	14.145	152	1207654	20.189	ng	99
50) Dimethylphthalate	13.880	163	1005157	20.271	ng	100
51) 2,6-Dinitrotoluene	13.992	165	212702	20.474	ng	98
52) Acenaphthene	14.486	154	765635	20.122	ng	99
53) 3-Nitroaniline	14.322	138	203884	20.306	ng	98
54) 2,4-Dinitrophenol	14.533	184	104048	19.806	ng	97
55) Dibenzofuran	14.822	168	1280124	20.125	ng	99
56) 4-Nitrophenol	14.639	139	180482	20.171	ng	98
57) 2,4-Dinitrotoluene	14.780	165	277724	20.039	ng	99
58) Fluorene	15.469	166	1019442	20.163	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.051	232	292142	19.603	ng	99
60) Diethylphthalate	15.245	149	968568	20.369	ng	100
61) 4-Chlorophenyl-phenyle...	15.463	204	535017	19.737	ng	96
62) 4-Nitroaniline	15.480	138	211717	20.390	ng	93
63) Azobenzene	15.757	77	832670	20.453	ng	99
65) 4,6-Dinitro-2-methylph...	15.545	198	164040	18.880	ng	94
66) n-Nitrosodiphenylamine	15.680	169	835415	20.246	ng	100
67) 4-Bromophenyl-phenylether	16.357	248	312714	19.537	ng	97
68) Hexachlorobenzene	16.474	284	361698	19.701	ng	96
69) Atrazine	16.627	200	195306	18.255	ng	98
70) Pentachlorophenol	16.821	266	263410	19.488	ng	99
71) Phenanthrene	17.210	178	1510233	19.979	ng	99
72) Anthracene	17.298	178	1481463	19.887	ng	100
73) Carbazole	17.568	167	1411278	20.114	ng	98
74) Di-n-butylphthalate	18.133	149	1638787	20.278	ng	99
75) Fluoranthene	19.221	202	1806131	19.433	ng	99
77) Benzidine	19.409	184	274181	15.426	ng	99
78) Pyrene	19.586	202	1868988	20.242	ng	99
80) Butylbenzylphthalate	20.486	149	719491	20.738	ng	98
81) Benzo(a)anthracene	21.386	228	1788846	20.090	ng	100
82) 3,3'-Dichlorobenzidine	21.309	252	658767	19.592	ng	99
83) Chrysene	21.445	228	1689336	19.957	ng	100
84) Bis(2-ethylhexyl)phtha...	21.309	149	1058190	20.934	ng	99
85) Di-n-octyl phthalate	22.445	149	1830449	20.699	ng	100
87) Indeno(1,2,3-cd)pyrene	27.785	276	2199957	19.848	ng	99
88) Benzo(b)fluoranthene	23.456	252	1818051	19.144	ng	99
89) Benzo(k)fluoranthene	23.521	252	1797007	19.527	ng	99
90) Benzo(a)pyrene	24.262	252	1629441	19.526	ng	99
91) Dibenzo(a,h)anthracene	27.844	278	1792231	19.631	ng	100
92) Benzo(g,h,i)perylene	28.844	276	1866337	20.005	ng	98

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

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