

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020725\
 Data File : BM049603.D
 Acq On : 07 Feb 2025 15:43
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Feb 07 16:35:25 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	191851	20.000	ng	0.00
21) Naphthalene-d8	10.622	136	735469	20.000	ng	0.00
39) Acenaphthene-d10	14.463	164	476309	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	903276	20.000	ng	0.00
76) Chrysene-d12	21.439	240	887934	20.000	ng	0.00
86) Perylene-d12	24.468	264	715919	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	971833	81.459	ng	0.00
7) Phenol-d6	6.981	99	1299547	83.143	ng	0.00
23) Nitrobenzene-d5	8.975	82	1277873	89.372	ng	0.00
42) 2,4,6-Tribromophenol	15.945	330	446090	79.040	ng	0.00
45) 2-Fluorobiphenyl	13.092	172	3195951	86.923	ng	0.00
79) Terphenyl-d14	19.827	244	5607445	96.435	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.293	88	208757	41.616	ng	99
3) Pyridine	3.693	79	569407	40.849	ng	100
4) n-Nitrosodimethylamine	3.599	42	277430	42.539	ng	92
6) Aniline	7.145	93	580828	39.208	ng	99
8) 2-Chlorophenol	7.387	128	481959	40.457	ng	96
9) Benzaldehyde	6.957	77	338401	40.889	ng	98
10) Phenol	7.010	94	622921	40.462	ng	99
11) bis(2-Chloroethyl)ether	7.245	93	518342	41.618	ng	96
12) 1,3-Dichlorobenzene	7.716	146	572239	41.384	ng	99
13) 1,4-Dichlorobenzene	7.857	146	582138	41.622	ng	99
14) 1,2-Dichlorobenzene	8.175	146	571131	42.290	ng	100
15) Benzyl Alcohol	8.057	79	420703	38.688	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.357	45	728137	43.428	ng	99
17) 2-Methylphenol	8.257	107	391275	41.566	ng	98
18) Hexachloroethane	8.910	117	215179	42.119	ng	98
19) n-Nitroso-di-n-propyla...	8.634	70	450106	43.976	ng	97
20) 3+4-Methylphenols	8.587	107	516749	41.584	ng	98
22) Acetophenone	8.639	105	833319	42.067	ng	# 97
24) Nitrobenzene	9.022	77	607023	43.696	ng	99
25) Isophorone	9.545	82	1029542	39.925	ng	98
26) 2-Nitrophenol	9.733	139	156680	35.341	ng	98
27) 2,4-Dimethylphenol	9.792	122	307484	37.712	ng	98
28) bis(2-Chloroethoxy)met...	10.033	93	688260	41.441	ng	97
29) 2,4-Dichlorophenol	10.263	162	460863	41.043	ng	98
30) 1,2,4-Trichlorobenzene	10.486	180	561616	40.824	ng	99
31) Naphthalene	10.669	128	1592814	41.064	ng	99
32) Benzoic acid	9.910	122	176299m	38.615	ng	
33) 4-Chloroaniline	10.769	127	571689	39.376	ng	99
34) Hexachlorobutadiene	10.969	225	340647	41.350	ng	99
35) Caprolactam	11.539	113	128576	37.534	ng	90
36) 4-Chloro-3-methylphenol	11.898	107	443383	39.595	ng	95
37) 2-Methylnaphthalene	12.286	142	1126789	41.259	ng	98
38) 1-Methylnaphthalene	12.504	142	1092548	41.117	ng	99
40) 1,2,4,5-Tetrachloroben...	12.651	216	643127	41.067	ng	99
41) Hexachlorocyclopentadiene	12.645	237	143229	35.827	ng	99
43) 2,4,6-Trichlorophenol	12.892	196	329641	38.574	ng	99

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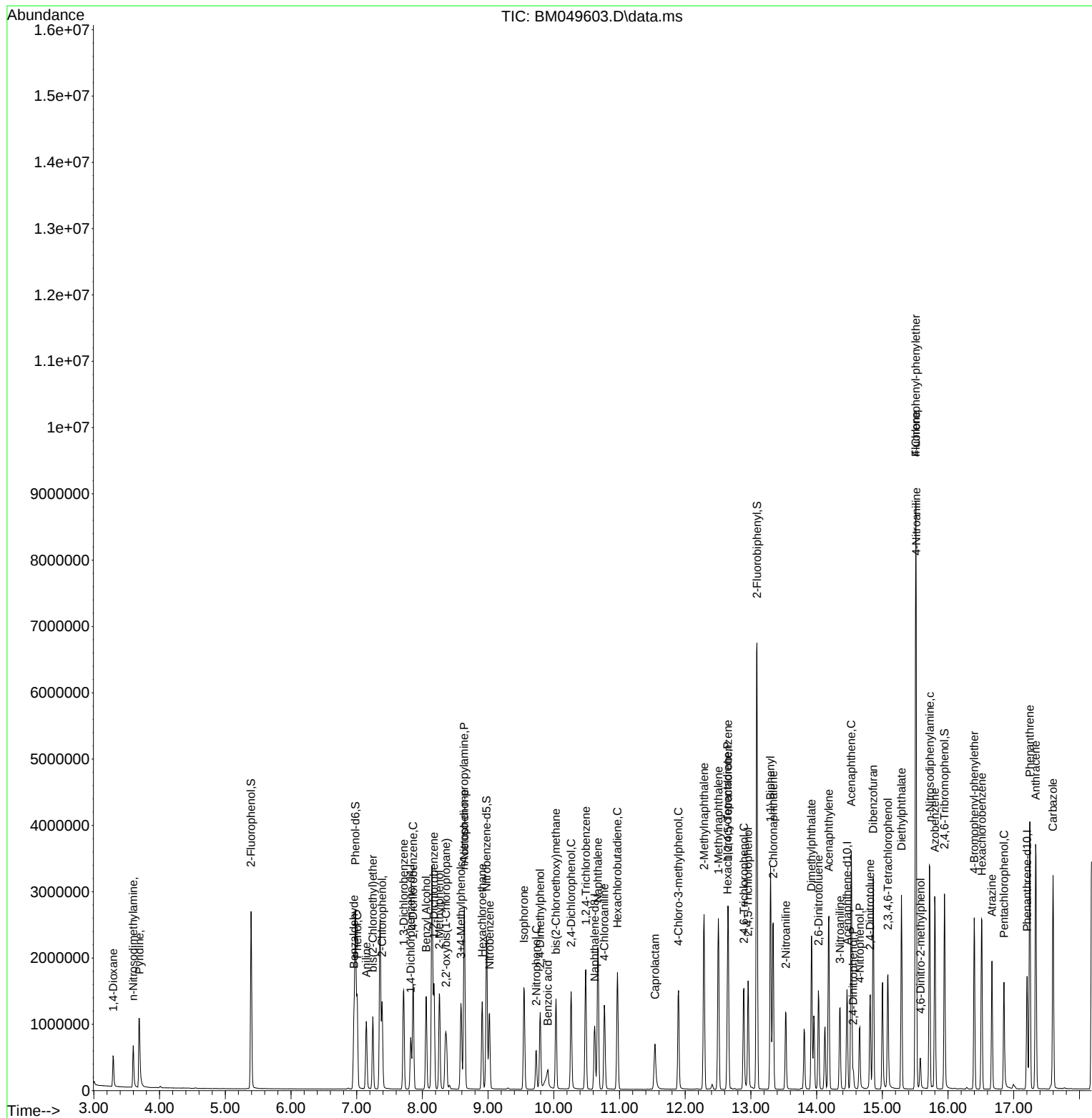
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.957	196	373626	39.785	ng	96
46) 1,1'-Biphenyl	13.298	154	1497729	41.219	ng	99
47) 2-Chloronaphthalene	13.339	162	1159896	41.259	ng	99
48) 2-Nitroaniline	13.533	65	281559	38.847	ng	97
49) Acenaphthylene	14.186	152	1702324	41.065	ng	99
50) Dimethylphthalate	13.921	163	1366535	39.950	ng	100
51) 2,6-Dinitrotoluene	14.027	165	258723	43.711	ng	94
52) Acenaphthene	14.527	154	1146562	41.669	ng	98
53) 3-Nitroaniline	14.357	138	265617	44.077	ng #	96
54) 2,4-Dinitrophenol	14.557	184	46156	35.883	ng	87
55) Dibenzofuran	14.863	168	1842163	42.046	ng	100
56) 4-Nitrophenol	14.657	139	197868	38.951	ng	97
57) 2,4-Dinitrotoluene	14.816	165	337426	42.109	ng	99
58) Fluorene	15.510	166	1696934	45.275	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.086	232	328177	41.362	ng	100
60) Diethylphthalate	15.292	149	1410288	41.025	ng	99
61) 4-Chlorophenyl-phenyle...	15.510	204	897185	44.974	ng	97
62) 4-Nitroaniline	15.515	138	324709	42.196	ng	99
63) Azobenzene	15.798	77	1573574	43.855	ng	97
65) 4,6-Dinitro-2-methylph...	15.580	198	79440	35.253	ng	91
66) n-Nitrosodiphenylamine	15.721	169	1178496	41.370	ng	99
67) 4-Bromophenyl-phenylether	16.398	248	429202	40.901	ng	100
68) Hexachlorobenzene	16.515	284	484650	40.239	ng	100
69) Atrazine	16.668	200	295969	34.952	ng	96
70) Pentachlorophenol	16.851	266	239189	34.300	ng	100
71) Phenanthrene	17.245	178	2156248	41.902	ng	100
72) Anthracene	17.333	178	2112998	41.470	ng	100
73) Carbazole	17.598	167	1818068	40.654	ng	99
74) Di-n-butylphthalate	18.186	149	2291870	40.388	ng	100
75) Fluoranthene	19.256	202	2501450	41.520	ng	99
77) Benzidine	19.439	184	251608	31.998	ng	96
78) Pyrene	19.621	202	2616645	39.357	ng	99
80) Butylbenzylphthalate	20.533	149	662322	32.014	ng	97
81) Benzo(a)anthracene	21.421	228	2441517	40.572	ng	99
82) 3,3'-Dichlorobenzidine	21.345	252	556530	34.368	ng	98
83) Chrysene	21.486	228	2313800	41.040	ng	100
84) Bis(2-ethylhexyl)phtha...	21.374	149	1282753	36.169	ng	97
85) Di-n-octyl phthalate	22.527	149	1706325	30.496	ng	97
87) Indeno(1,2,3-cd)pyrene	27.891	276	1587030	40.591	ng	99
88) Benzo(b)fluoranthene	23.515	252	2027116	41.159	ng	99
89) Benzo(k)fluoranthene	23.580	252	1939592	40.547	ng	100
90) Benzo(a)pyrene	24.327	252	1574267	40.335	ng	100
91) Dibenzo(a,h)anthracene	27.956	278	1249408	40.433	ng	97
92) Benzo(g,h,i)perylene	28.956	276	1430421	41.439	ng	98

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

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