

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042825\
 Data File : BM050031.D
 Acq On : 28 Apr 2025 17:04
 Operator : RC/JU
 Sample : SSTDICC080
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC080

Quant Time: Apr 28 17:59:20 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM042825.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 28 17:38:08 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.769	152	251380	20.000	ng	0.00
21) Naphthalene-d8	10.563	136	931435	20.000	ng	0.00
39) Acenaphthene-d10	14.410	164	613490	20.000	ng	0.00
64) Phenanthrene-d10	17.156	188	1212251	20.000	ng	0.00
76) Chrysene-d12	21.397	240	1197780	20.000	ng	0.00
86) Perylene-d12	24.397	264	1141905	20.000	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.357	112	2297313	160.027	ng	0.00
7) Phenol-d6	6.957	99	2954036	163.719	ng	0.00
23) Nitrobenzene-d5	8.933	82	2984192	157.875	ng	0.00
42) 2,4,6-Tribromophenol	15.909	330	1586083	174.847	ng	0.00
45) 2-Fluorobiphenyl	13.033	172	8283812	159.736	ng	0.00
79) Terphenyl-d14	19.780	244	11877110	146.730	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	3.251	88	475912	77.317	ng	99
3) Pyridine	3.657	79	1267903	80.171	ng	99
4) n-Nitrosodimethylamine	3.563	42	492501	79.400	ng	99
6) Aniline	7.098	93	1845033	80.979	ng	99
8) 2-Chlorophenol	7.339	128	1241711	79.998	ng	100
9) Benzaldehyde	6.910	77	898993	74.544	ng	99
10) Phenol	6.981	94	1491194	81.638	ng	98
11) bis(2-Chloroethyl)ether	7.192	93	1224517	80.283	ng	99
12) 1,3-Dichlorobenzene	7.657	146	1460820	77.912	ng	100
13) 1,4-Dichlorobenzene	7.804	146	1465881	77.822	ng	99
14) 1,2-Dichlorobenzene	8.122	146	1412270	77.618	ng	99
15) Benzyl Alcohol	8.016	79	1055328	84.339	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.292	45	1440484	77.694	ng	100
17) 2-Methylphenol	8.228	107	980375	81.473	ng	99
18) Hexachloroethane	8.845	117	513126	76.632	ng	98
19) n-Nitroso-di-n-propyla...	8.581	70	933468	80.665	ng	99
20) 3+4-Methylphenols	8.557	107	1348190	82.965	ng	99
22) Acetophenone	8.592	105	1815657	78.406	ng	# 99
24) Nitrobenzene	8.975	77	1289068	78.871	ng	99
25) Isophorone	9.504	82	2559279	79.430	ng	100
26) 2-Nitrophenol	9.680	139	710174	85.053	ng	99
27) 2,4-Dimethylphenol	9.751	122	1136040	82.043	ng	99
28) bis(2-Chloroethoxy)met...	9.975	93	1581985	79.437	ng	99
29) 2,4-Dichlorophenol	10.222	162	1275806	82.243	ng	99
30) 1,2,4-Trichlorobenzene	10.427	180	1485344	78.590	ng	99
31) Naphthalene	10.610	128	3686066	78.143	ng	100
32) Benzoic acid	9.951	122	800660	90.660	ng	97
33) 4-Chloroaniline	10.727	127	1575909	81.508	ng	99
34) Hexachlorobutadiene	10.898	225	959905	80.012	ng	99
35) Caprolactam	11.539	113	371164	82.060	ng	98
36) 4-Chloro-3-methylphenol	11.869	107	1152624	80.424	ng	99
37) 2-Methylnaphthalene	12.227	142	2525532	79.858	ng	99
38) 1-Methylnaphthalene	12.451	142	2645483	79.045	ng	99
40) 1,2,4,5-Tetrachloroben...	12.604	216	1799698	84.742	ng	99
41) Hexachlorocyclopentadiene	12.580	237	1185418	91.336	ng	99
43) 2,4,6-Trichlorophenol	12.851	196	1151641	85.496	ng	99

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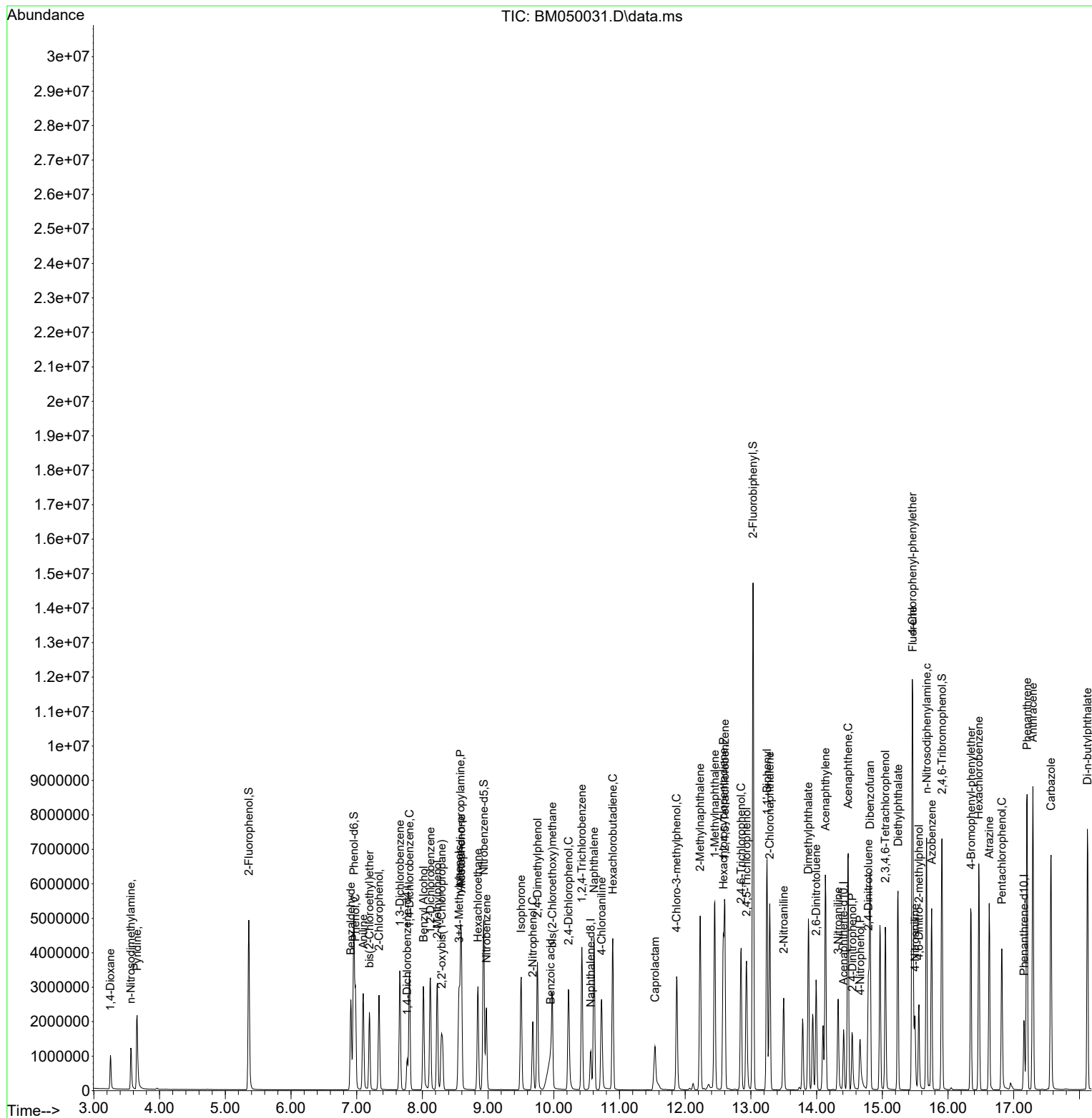
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.933	196	1220306	84.011	ng	98
46) 1,1'-Biphenyl	13.245	154	3613893	79.246	ng	99
47) 2-Chloronaphthalene	13.286	162	2842514	78.700	ng	99
48) 2-Nitroaniline	13.498	65	701256	82.725	ng	99
49) Acenaphthylene	14.133	152	4526874	79.492	ng	99
50) Dimethylphthalate	13.874	163	3507333	78.228	ng	99
51) 2,6-Dinitrotoluene	13.992	165	758296	81.689	ng	99
52) Acenaphthene	14.480	154	2618967	79.452	ng	100
53) 3-Nitroaniline	14.327	138	740202	83.521	ng	100
54) 2,4-Dinitrophenol	14.545	184	469472	79.096	ng	98
55) Dibenzofuran	14.815	168	4325783	78.875	ng	99
56) 4-Nitrophenol	14.662	139	566943	78.710	ng	99
57) 2,4-Dinitrotoluene	14.786	165	1080118	84.207	ng	96
58) Fluorene	15.462	166	3565506	79.426	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.045	232	1057707	84.123	ng	96
60) Diethylphthalate	15.239	149	3257954	77.118	ng	97
61) 4-Chlorophenyl-phenyle...	15.457	204	2051105	83.186	ng	96
62) 4-Nitroaniline	15.498	138	724549	84.678	ng	99
63) Azobenzene	15.751	77	2723186	76.914	ng	95
65) 4,6-Dinitro-2-methylph...	15.557	198	665798	87.741	ng	96
66) n-Nitrosodiphenylamine	15.674	169	2957293	79.443	ng	99
67) 4-Bromophenyl-phenylether	16.351	248	1225531	83.600	ng	94
68) Hexachlorobenzene	16.468	284	1412768	83.606	ng	95
69) Atrazine	16.627	200	1116818	83.174	ng	100
70) Pentachlorophenol	16.821	266	843195	88.648	ng	99
71) Phenanthrene	17.203	178	5487968	80.266	ng	99
72) Anthracene	17.292	178	5597195	80.893	ng	99
73) Carbazole	17.568	167	4845637	80.705	ng	99
74) Di-n-butylphthalate	18.121	149	5649166	79.007	ng	99
75) Fluoranthene	19.221	202	6841999	85.234	ng	99
77) Benzidine	19.403	184	3598340	84.575	ng	100
78) Pyrene	19.580	202	6959279	82.738	ng	99
80) Butylbenzylphthalate	20.474	149	2512705	79.760	ng	97
81) Benzo(a)anthracene	21.380	228	6747718	81.808	ng	99
82) 3,3'-Dichlorobenzidine	21.303	252	2782107	88.111	ng	99
83) Chrysene	21.444	228	6207632	81.316	ng	98
84) Bis(2-ethylhexyl)phtha...	21.297	149	3665836	77.905	ng	# 96
85) Di-n-octyl phthalate	22.427	149	6109632	78.732	ng	99
87) Indeno(1,2,3-cd)pyrene	27.803	276	7285836	85.629	ng	98
88) Benzo(b)fluoranthene	23.456	252	6347901	85.480	ng	99
89) Benzo(k)fluoranthene	23.521	252	6265924	83.396	ng	99
90) Benzo(a)pyrene	24.262	252	5925652	85.195	ng	99
91) Dibenzo(a,h)anthracene	27.850	278	5981478	86.239	ng	99
92) Benzo(g,h,i)perylene	28.862	276	5537637	83.395	ng	99

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

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