

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM020625\
 Data File : BM049592.D
 Acq On : 06 Feb 2025 23:34
 Operator : RC/JU
 Sample : Q1168-07
 Misc : LOD-MDL-SOIL 4PPM
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 LOD-MDL-WATER-01-QT1-2025

Quant Time: Feb 07 01:33:26 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

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 02/07/2025
 Supervised By :mohammad ahmed 02/07/2025

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	212080	20.000	ng	0.00
21) Naphthalene-d8	10.622	136	769359	20.000	ng	0.00
39) Acenaphthene-d10	14.463	164	471787	20.000	ng	0.00
64) Phenanthrene-d10	17.204	188	879190	20.000	ng	0.00
76) Chrysene-d12	21.439	240	727368	20.000	ng	0.00
86) Perylene-d12	24.468	264	613590	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.393	112	1727150	130.960	ng	0.00
7) Phenol-d6	6.981	99	2221875	128.593	ng	0.00
23) Nitrobenzene-d5	8.975	82	1473900	98.541	ng	0.00
42) 2,4,6-Tribromophenol	15.945	330	711385	127.254	ng	0.00
45) 2-Fluorobiphenyl	13.086	172	3593389	98.669	ng	0.00
79) Terphenyl-d14	19.827	244	6203797	130.243	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.299	88	25940	4.678	ng	98
3) Pyridine	3.705	79	56213	3.648	ng	# 92
4) n-Nitrosodimethylamine	3.605	42	29307	4.065	ng	# 83
6) Aniline	7.146	93	69275	4.230	ng	99
8) 2-Chlorophenol	7.387	128	54692	4.153	ng	99
9) Benzaldehyde	6.957	77	60846	6.651	ng	99
10) Phenol	7.005	94	70233	4.127	ng	90
11) bis(2-Chloroethyl)ether	7.246	93	58899	4.278	ng	95
12) 1,3-Dichlorobenzene	7.716	146	64732	4.235	ng	97
13) 1,4-Dichlorobenzene	7.857	146	65012	4.205	ng	98
14) 1,2-Dichlorobenzene	8.175	146	63475	4.252	ng	97
15) Benzyl Alcohol	8.057	79	40948	3.406	ng	90
16) 2,2'-oxybis(1-Chloropr...	8.346	45	85194	4.597	ng	98
17) 2-Methylphenol	8.257	107	38387	3.689	ng	95
18) Hexachloroethane	8.910	117	23274	4.121	ng	95
19) n-Nitroso-di-n-propyla...	8.628	70	41889	3.702	ng	96
20) 3+4-Methylphenols	8.581	107	45142	3.286	ng	96
22) Acetophenone	8.640	105	85956	4.148	ng	96
24) Nitrobenzene	9.016	77	71298	4.906	ng	96
25) Isophorone	9.546	82	105826	3.923	ng	99
26) 2-Nitrophenol	9.734	139	9802	4.313	ng	94
27) 2,4-Dimethylphenol	9.793	122	23061m	2.704	ng	
28) bis(2-Chloroethoxy)met...	10.034	93	67459	3.883	ng	98
29) 2,4-Dichlorophenol	10.263	162	36276	3.088	ng	96
30) 1,2,4-Trichlorobenzene	10.481	180	59554	4.138	ng	100
31) Naphthalene	10.669	128	159901	3.941	ng	96
32) Benzoic acid	9.828	122	6724m	6.543	ng	
33) 4-Chloroaniline	10.769	127	57029	3.755	ng	98
34) Hexachlorobutadiene	10.969	225	33216	3.854	ng	95
35) Caprolactam	11.510	113	8873m	2.476	ng	
36) 4-Chloro-3-methylphenol	11.892	107	33120	2.827	ng	98
37) 2-Methylnaphthalene	12.281	142	104584	3.661	ng	96
38) 1-Methylnaphthalene	12.504	142	97604	3.511	ng	100
40) 1,2,4,5-Tetrachloroben...	12.651	216	59699	3.849	ng	99
41) Hexachlorocyclopentadiene	12.645	237	10571	2.670	ng	97
43) 2,4,6-Trichlorophenol	12.886	196	20293	2.397	ng	95

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44) 2,4,5-Trichlorophenol	12.957	196	25693	2.762	ng	94
46) 1,1'-Biphenyl	13.298	154	143498	3.987	ng	98
47) 2-Chloronaphthalene	13.334	162	109616	3.937	ng	98
48) 2-Nitroaniline	13.528	65	17581	6.316	ng	90
49) Acenaphthylene	14.181	152	163882	3.991	ng	99
50) Dimethylphthalate	13.916	163	121759	3.594	ng	99
51) 2,6-Dinitrotoluene	14.022	165	17045	2.907	ng	# 80
52) Acenaphthene	14.522	154	95525	3.505	ng	95
53) 3-Nitroaniline	14.351	138	17605	2.949	ng	# 91
54) 2,4-Dinitrophenol	14.551	184	2293	4.618	ng	89
55) Dibenzofuran	14.857	168	160178	3.691	ng	98
57) 2,4-Dinitrotoluene	14.810	165	18507	5.159	ng	# 86
58) Fluorene	15.510	166	124124	3.343	ng	100
59) 2,3,4,6-Tetrachlorophenol	15.080	232	19760	2.514	ng	95
60) Diethylphthalate	15.286	149	112648	3.308	ng	97
61) 4-Chlorophenyl-phenyle...	15.504	204	64033	3.241	ng	91
62) 4-Nitroaniline	15.510	138	16638	6.176	ng	94
63) Azobenzene	15.798	77	132354	3.724	ng	97
65) 4,6-Dinitro-2-methylph...	15.575	198	3850	6.064	ng	91
66) n-Nitrosodiphenylamine	15.716	169	98164	3.540	ng	100
67) 4-Bromophenyl-phenylether	16.398	248	35747	3.500	ng	98
68) Hexachlorobenzene	16.510	284	42538	3.629	ng	96
69) Atrazine	16.669	200	27015	3.278	ng	94
70) Pentachlorophenol	16.845	266	12127m	9.098	ng	
71) Phenanthrene	17.245	178	188660	3.767	ng	98
72) Anthracene	17.333	178	172672	3.482	ng	99
73) Carbazole	17.598	167	157819	3.626	ng	100
74) Di-n-butylphthalate	18.186	149	144979	2.625	ng	99
75) Fluoranthene	19.257	202	188618	3.217	ng	99
77) Benzidine	19.439	184	16771	2.604	ng	# 95
78) Pyrene	19.615	202	202213	3.713	ng	99
81) Benzo(a)anthracene	21.421	228	180925	3.670	ng	98
82) 3,3'-Dichlorobenzidine	21.345	252	28388	2.140	ng	# 90
83) Chrysene	21.480	228	172705	3.739	ng	99
87) Indeno(1,2,3-cd)pyrene	27.897	276	126862m	3.786	ng	
88) Benzo(b)fluoranthene	23.509	252	136135	3.225	ng	98
89) Benzo(k)fluoranthene	23.574	252	143400	3.498	ng	98
90) Benzo(a)pyrene	24.327	252	112794	3.372	ng	99
91) Dibenzo(a,h)anthracene	27.962	278	96708m	3.652	ng	
92) Benzo(g,h,i)perylene	28.938	276	109011m	3.685	ng	

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

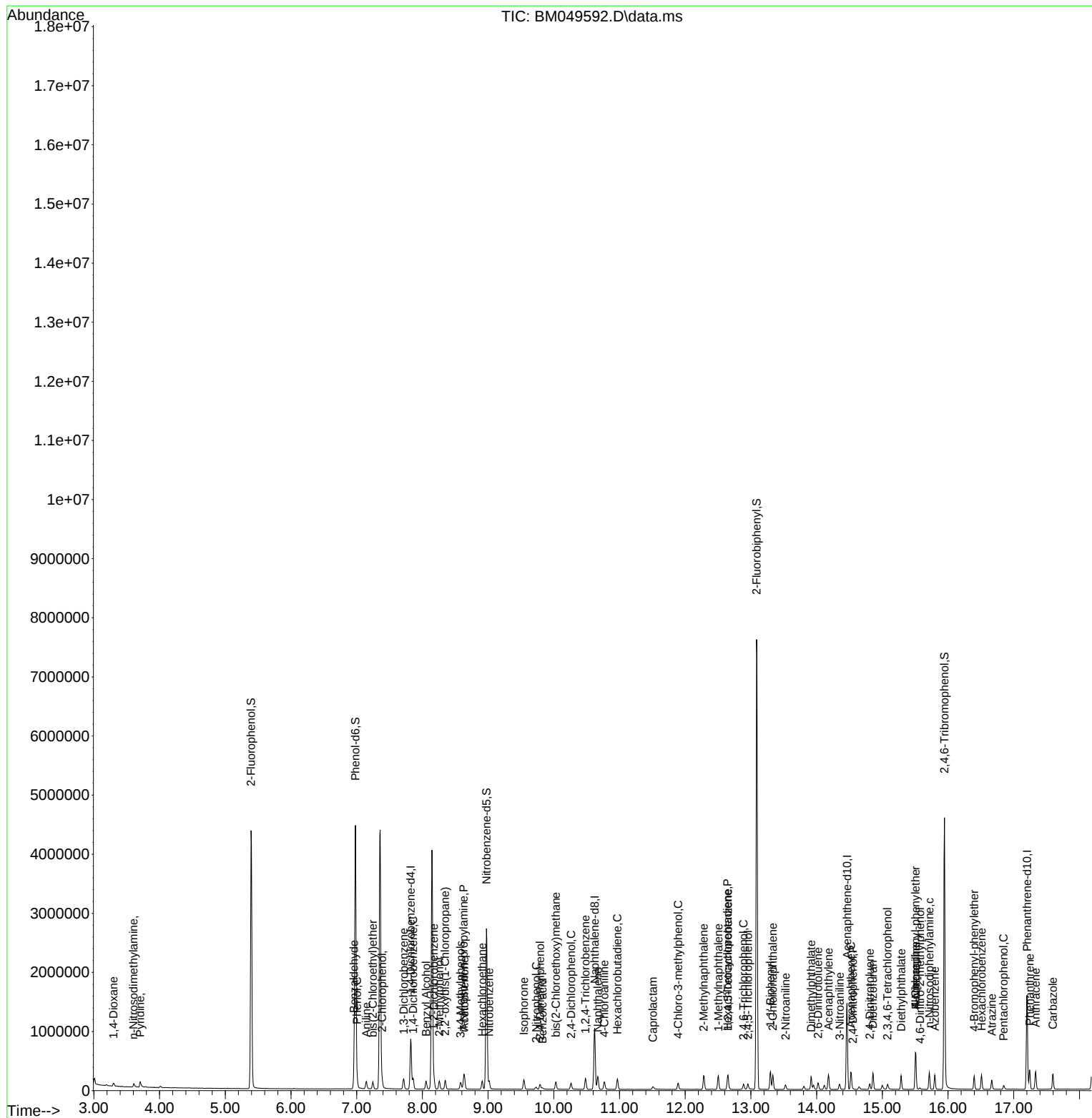
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