

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082224\
 Data File : BM047308.D
 Acq On : 22 Aug 2024 09:14
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Quant Time: Aug 22 10:59:43 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Aug 11 23:47:58 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.351	152	315372	20.000	ng	-0.02
21) Naphthalene-d8	10.110	136	1211803	20.000	ng	-0.02
39) Acenaphthene-d10	14.016	164	822573	20.000	ng	-0.02
64) Phenanthrene-d10	16.780	188	1768064	20.000	ng	-0.01
76) Chrysene-d12	21.021	240	1535454	20.000	ng	-0.01
86) Perylene-d12	23.709	264	1629252	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.010	112	1303996	73.882	ng	-0.01
7) Phenol-d6	6.569	99	1761491	72.430	ng	-0.02
23) Nitrobenzene-d5	8.504	82	1721029	71.513	ng	-0.02
42) 2,4,6-Tribromophenol	15.527	330	800401	83.918	ng	-0.01
45) 2-Fluorobiphenyl	12.622	172	4120450	77.269	ng	-0.02
79) Terphenyl-d14	19.445	244	6380197	89.037	ng	-0.02
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.028	88	263249	35.998	ng	96
3) Pyridine	3.399	79	734306	34.243	ng	98
4) n-Nitrosodimethylamine	3.322	42	306437	32.708	ng	96
6) Aniline	6.704	93	809995	35.238	ng	98
8) 2-Chlorophenol	6.928	128	714965	37.141	ng	99
9) Benzaldehyde	6.516	77	523783	41.837	ng	99
10) Phenol	6.593	94	863321	35.576	ng	98
11) bis(2-Chloroethyl)ether	6.804	93	742990	35.752	ng	97
12) 1,3-Dichlorobenzene	7.240	146	879161	38.433	ng	97
13) 1,4-Dichlorobenzene	7.387	146	886468	38.275	ng	98
14) 1,2-Dichlorobenzene	7.698	146	825943	37.796	ng	98
15) Benzyl Alcohol	7.604	79	619139	35.771	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.892	45	880322	33.012	ng	98
17) 2-Methylphenol	7.816	107	598616	36.976	ng	99
18) Hexachloroethane	8.404	117	331304	36.929	ng	94
19) n-Nitroso-di-n-propyla...	8.169	70	561893	34.301	ng	96
20) 3+4-Methylphenols	8.145	107	818840	36.179	ng	100
22) Acetophenone	8.175	105	1094687	37.424	ng	98
24) Nitrobenzene	8.545	77	884972	36.519	ng	97
25) Isophorone	9.069	82	1591106	37.971	ng	98
26) 2-Nitrophenol	9.245	139	449274	41.863	ng	94
27) 2,4-Dimethylphenol	9.328	122	496910	39.527	ng	99
28) bis(2-Chloroethoxy)met...	9.551	93	993074	37.664	ng	100
29) 2,4-Dichlorophenol	9.775	162	782273	41.938	ng	99
30) 1,2,4-Trichlorobenzene	9.975	180	881137	41.742	ng	100
31) Naphthalene	10.157	128	2303393	38.156	ng	99
32) Benzoic acid	9.522	122	603943	41.350	ng	99
33) 4-Chloroaniline	10.286	127	873816	37.641	ng	99
34) Hexachlorobutadiene	10.439	225	554015	44.840	ng	100
35) Caprolactam	11.104	113	259311	40.031	ng	91
36) 4-Chloro-3-methylphenol	11.439	107	785639	39.548	ng	100
37) 2-Methylnaphthalene	11.786	142	1687918	39.270	ng	98
38) 1-Methylnaphthalene	12.010	142	1643957	38.700	ng	99
40) 1,2,4,5-Tetrachloroben...	12.169	216	957174	42.662	ng	98
41) Hexachlorocyclopentadiene	12.139	237	284615	36.329	ng	98
43) 2,4,6-Trichlorophenol	12.427	196	665686	41.744	ng	100

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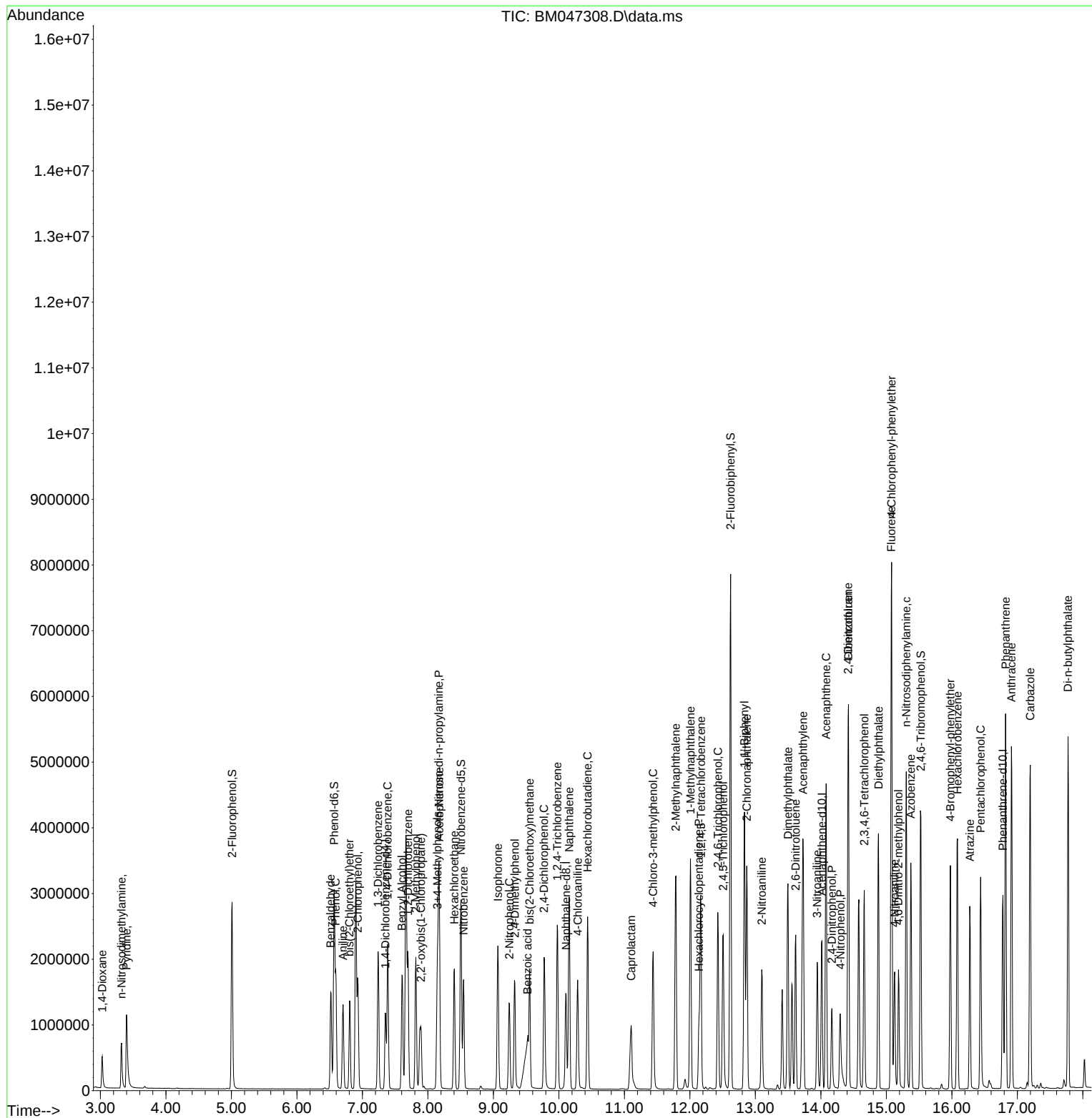
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.510	196	738243	41.695	ng	96
46) 1,1'-Biphenyl	12.833	154	2220122	37.912	ng	98
47) 2-Chloronaphthalene	12.869	162	1733753	37.879	ng	99
48) 2-Nitroaniline	13.098	65	519350	36.219	ng	96
49) Acenaphthylene	13.727	152	2561019	37.959	ng	100
50) Dimethylphthalate	13.498	163	2232553	38.861	ng	100
51) 2,6-Dinitrotoluene	13.616	165	536890	39.978	ng	88
52) Acenaphthene	14.080	154	1620780	37.872	ng	100
53) 3-Nitroaniline	13.945	138	510904	37.966	ng	100
54) 2,4-Dinitrophenol	14.169	184	303462	37.702	ng	93
55) Dibenzofuran	14.421	168	2590656	38.323	ng	99
56) 4-Nitrophenol	14.298	139	415593	35.815	ng	98
57) 2,4-Dinitrotoluene	14.416	165	708560	39.529	ng	# 97
58) Fluorene	15.074	166	2133336	38.248	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.663	232	609310	41.872	ng	93
60) Diethylphthalate	14.880	149	2237417	38.138	ng	98
61) 4-Chlorophenyl-phenyle...	15.080	204	1071908	40.367	ng	95
62) 4-Nitroaniline	15.127	138	500966	37.788	ng	95
63) Azobenzene	15.374	77	1953128	35.112	ng	95
65) 4,6-Dinitro-2-methylph...	15.186	198	425285	41.084	ng	92
66) n-Nitrosodiphenylamine	15.304	169	1831258	38.026	ng	99
67) 4-Bromophenyl-phenylether	15.980	248	705879	41.431	ng	96
68) Hexachlorobenzene	16.086	284	821372	40.254	ng	94
69) Atrazine	16.274	200	511524	35.298	ng	99
70) Pentachlorophenol	16.439	266	580674	40.597	ng	98
71) Phenanthrene	16.821	178	3343698	37.683	ng	100
72) Anthracene	16.910	178	3284636	38.044	ng	100
73) Carbazole	17.198	167	3281091	37.220	ng	99
74) Di-n-butylphthalate	17.774	149	4010258	39.097	ng	99
75) Fluoranthene	18.856	202	4234649	39.236	ng	99
77) Benzidine	19.062	184	1551050	59.567	ng	100
78) Pyrene	19.221	202	4349870	39.676	ng	100
80) Butylbenzylphthalate	20.162	149	1852458	41.403	ng	93
81) Benzo(a)anthracene	21.003	228	3891993	38.182	ng	100
82) 3,3'-Dichlorobenzidine	20.945	252	1389102	43.555	ng	99
83) Chrysene	21.062	228	3637684	37.614	ng	100
84) Bis(2-ethylhexyl)phtha...	20.956	149	2572281	40.539	ng	98
85) Di-n-octyl phthalate	21.986	149	4548357	42.173	ng	98
87) Indeno(1,2,3-cd)pyrene	26.703	276	3627893	34.449	ng	98
88) Benzo(b)fluoranthene	22.874	252	3826072	39.544	ng	98
89) Benzo(k)fluoranthene	22.927	252	3657262	40.018	ng	99
90) Benzo(a)pyrene	23.586	252	3250485	38.781	ng	98
91) Dibenzo(a,h)anthracene	26.750	278	2934810	34.436	ng	98
92) Benzo(g,h,i)perylene	27.638	276	2982909	33.711	ng	98

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

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