

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060625\
 Data File : BM050213.D
 Acq On : 06 Jun 2025 14:06
 Operator : RC/JU
 Sample : Q2194-02MS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 COMP-12MS

Quant Time: Jun 06 14:37:50 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 05 16:20:25 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	348615	20.000	ng	0.00
21) Naphthalene-d8	10.575	136	1377434	20.000	ng	0.00
39) Acenaphthene-d10	14.416	164	768114	20.000	ng	0.00
64) Phenanthrene-d10	17.151	188	1293379	20.000	ng	0.00
76) Chrysene-d12	21.380	240	1106417	20.000	ng	0.00
86) Perylene-d12	24.368	264	1212630	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.375	112	2485049	118.911	ng	0.00
7) Phenol-d6	6.957	99	2994886	108.862	ng	0.00
23) Nitrobenzene-d5	8.939	82	2093769	79.055	ng	0.00
42) 2,4,6-Tribromophenol	15.898	330	1069797	122.448	ng	0.00
45) 2-Fluorobiphenyl	13.045	172	4313578	76.030	ng	0.00
79) Terphenyl-d14	19.780	244	4512257	77.284	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.304	88	309499	33.786	ng	# 78
3) Pyridine	3.693	79	837934	35.322	ng	100
4) n-Nitrosodimethylamine	3.599	42	200375	40.847	ng	95
6) Aniline	7.116	93	795951	21.886	ng	98
8) 2-Chlorophenol	7.351	128	994911	43.281	ng	99
9) Benzaldehyde	6.928	77	409612	23.418	ng	99
10) Phenol	6.981	94	1102221	37.866	ng	99
11) bis(2-Chloroethyl)ether	7.216	93	980654	41.885	ng	99
12) 1,3-Dichlorobenzene	7.681	146	775911	30.193	ng	99
13) 1,4-Dichlorobenzene	7.822	146	818226	30.602	ng	99
14) 1,2-Dichlorobenzene	8.139	146	807074	31.663	ng	99
15) Benzyl Alcohol	8.022	79	858483	43.752	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.316	45	657839	40.447	ng	98
17) 2-Methylphenol	8.228	107	812312	42.295	ng	100
18) Hexachloroethane	8.869	117	288971	29.843	ng	99
19) n-Nitroso-di-n-propyla...	8.592	70	713472	42.380	ng	99
20) 3+4-Methylphenols	8.551	107	1079323	42.004	ng	99
22) Acetophenone	8.604	105	1456030	42.311	ng	99
24) Nitrobenzene	8.981	77	1045437	43.402	ng	98
25) Isophorone	9.510	82	2007832	43.923	ng	100
26) 2-Nitrophenol	9.686	139	475401	45.722	ng	98
27) 2,4-Dimethylphenol	9.751	122	934333	43.734	ng	98
28) bis(2-Chloroethoxy)met...	9.992	93	1306450	43.748	ng	99
29) 2,4-Dichlorophenol	10.222	162	887017	44.867	ng	99
30) 1,2,4-Trichlorobenzene	10.439	180	830875	37.767	ng	99
31) Naphthalene	10.628	128	2714278	38.113	ng	100
32) Benzoic acid	9.845	122	281160	20.941	ng	98
33) 4-Chloroaniline	10.728	127	208080	6.853	ng	99
34) Hexachlorobutadiene	10.922	225	451045	34.468	ng	99
35) Caprolactam	11.498	113	211120	33.684	ng	97
36) 4-Chloro-3-methylphenol	11.851	107	909622	43.119	ng	100
37) 2-Methylnaphthalene	12.239	142	1787634	41.609	ng	98
38) 1-Methylnaphthalene	12.457	142	1857710	40.739	ng	99
40) 1,2,4,5-Tetrachloroben...	12.610	216	968442	43.663	ng	99
41) Hexachlorocyclopentadiene	12.598	237	1125098	83.526	ng	98
43) 2,4,6-Trichlorophenol	12.845	196	676045	46.349	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060625\
 Data File : BM050213.D
 Acq On : 06 Jun 2025 14:06
 Operator : RC/JU
 Sample : Q2194-02MS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 COMP-12MS

Quant Time: Jun 06 14:37:50 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 05 16:20:25 2025
 Response via : Initial Calibration

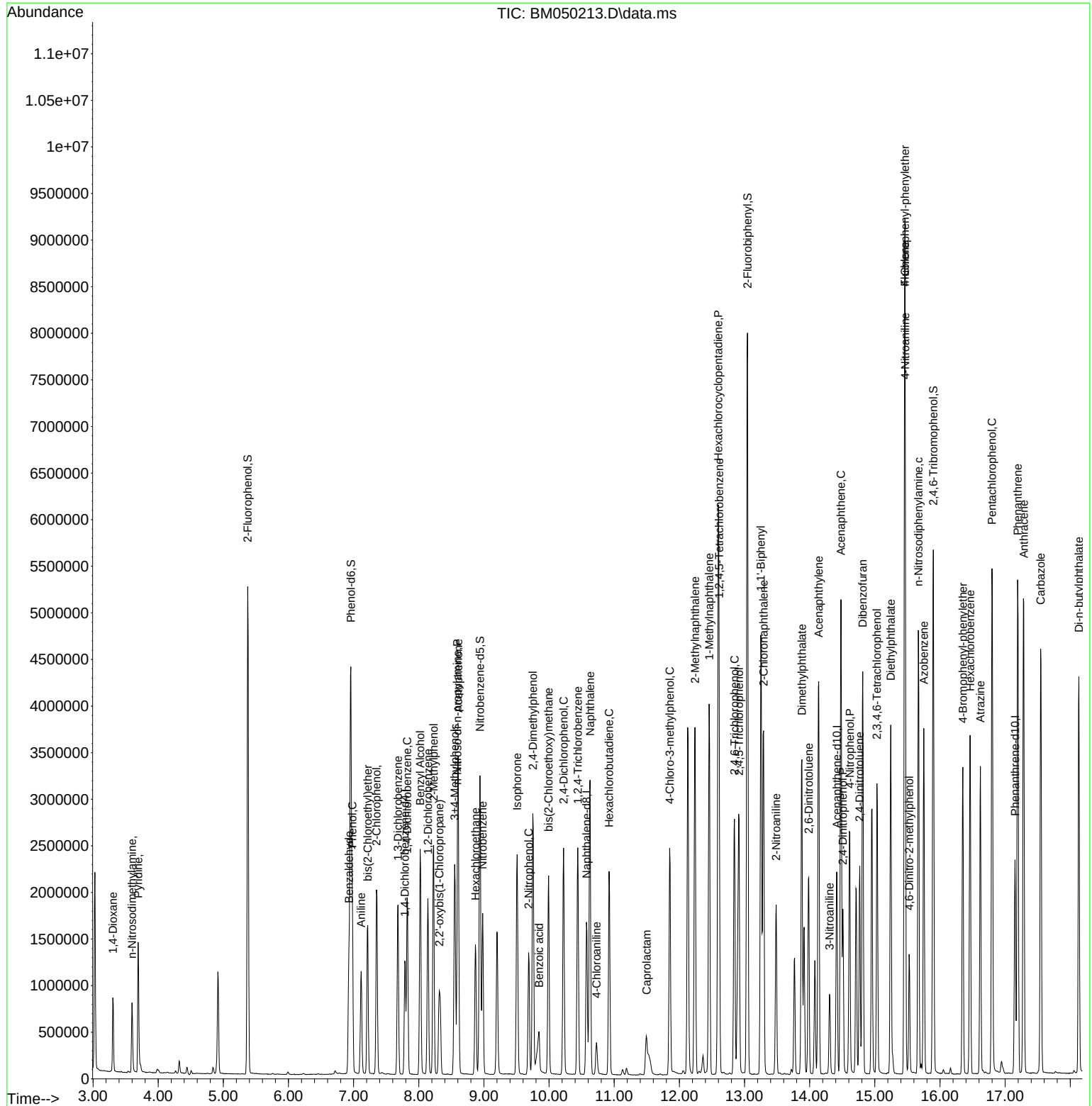
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.916	196	729356	45.543	ng	98
46) 1,1'-Biphenyl	13.251	154	2530450	43.996	ng	98
47) 2-Chloronaphthalene	13.292	162	1935364	43.283	ng	100
48) 2-Nitroaniline	13.486	65	477197	48.495	ng	97
49) Acenaphthylene	14.139	152	3153424	43.602	ng	100
50) Dimethylphthalate	13.880	163	2278169	43.855	ng	100
51) 2,6-Dinitrotoluene	13.986	165	488399	46.620	ng	98
52) Acenaphthene	14.480	154	2043709	45.178	ng	99
53) 3-Nitroaniline	14.304	138	232659	19.927	ng	96
54) 2,4-Dinitrophenol	14.510	184	379194	59.986	ng	98
55) Dibenzofuran	14.816	168	2842304	42.948	ng	99
56) 4-Nitrophenol	14.616	139	779304	78.421	ng	97
57) 2,4-Dinitrotoluene	14.768	165	648380	46.918	ng	99
58) Fluorene	15.463	166	2148065	42.377	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.033	232	599098	43.239	ng	92
60) Diethylphthalate	15.245	149	2186884	42.429	ng	99
61) 4-Chlorophenyl-phenyle...	15.463	204	1049834	42.854	ng	97
62) 4-Nitroaniline	15.468	138	426152	38.942	ng	98
63) Azobenzene	15.751	77	2199804	42.700	ng	99
65) 4,6-Dinitro-2-methylph...	15.527	198	287886	40.753	ng	97
66) n-Nitrosodiphenylamine	15.668	169	1858818	45.740	ng	99
67) 4-Bromophenyl-phenylether	16.351	248	651215	47.215	ng	98
68) Hexachlorobenzene	16.462	284	746011	46.295	ng	99
69) Atrazine	16.621	200	607699	46.931	ng	100
70) Pentachlorophenol	16.798	266	960880	91.715	ng	99
71) Phenanthrene	17.192	178	3258540	44.102	ng	99
72) Anthracene	17.286	178	3285651	44.451	ng	99
73) Carbazole	17.545	167	2919116	43.350	ng	99
74) Di-n-butylphthalate	18.133	149	3333066	45.491	ng	100
75) Fluoranthene	19.203	202	3339548	42.860	ng	99
77) Benzidine	19.386	184	595875	19.152	ng	98
78) Pyrene	19.568	202	3471185	46.548	ng	99
80) Butylbenzylphthalate	20.480	149	1237003	49.729	ng	99
81) Benzo(a)anthracene	21.362	228	3166143	45.210	ng	100
82) 3,3'-Dichlorobenzidine	21.286	252	469675	21.992	ng	99
83) Chrysene	21.427	228	3012726	45.226	ng	98
84) Bis(2-ethylhexyl)phtha...	21.315	149	1858055	48.493	ng	99
85) Di-n-octyl phthalate	22.456	149	2955416	51.980	ng	100
87) Indeno(1,2,3-cd)pyrene	27.738	276	3987489	51.070	ng	# 92
88) Benzo(b)fluoranthene	23.433	252	3197636	45.355	ng	99
89) Benzo(k)fluoranthene	23.491	252	3221599	44.761	ng	100
90) Benzo(a)pyrene	24.233	252	3082761	46.506	ng	100
91) Dibenzo(a,h)anthracene	27.809	278	3199032	50.476	ng	99
92) Benzo(g,h,i)perylene	28.791	276	3296635	51.971	ng	100

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060625\
Data File : BM050213.D
Acq On : 06 Jun 2025 14:06
Operator : RC/JU
Sample : Q2194-02MS
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COMP-12MS

Quant Time: Jun 06 14:37:50 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 05 16:20:25 2025
Response via : Initial Calibration



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060625\
Data File : BM050213.D
Acq On : 06 Jun 2025 14:06
Operator : RC/JU
Sample : Q2194-02MS
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COMP-12MS

Quant Time: Jun 06 14:37:50 2025
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM060525.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 05 16:20:25 2025
Response via : Initial Calibration

