

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060525\
 Data File : BM050196.D
 Acq On : 05 Jun 2025 10:38
 Operator : RC/JU
 Sample : SSTDICC010
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC010

Manual Integrations
 APPROVED

Reviewed By :Rahul Chavli 06/05/2025
 Supervised By :Jagrut Upadhyay 06/06/2025

Quant Time: Jun 05 14:58:04 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 14:55:18 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	257321	20.000	ng	0.00
21) Naphthalene-d8	10.575	136	1031213	20.000	ng	0.00
39) Acenaphthene-d10	14.416	164	597701	20.000	ng	0.00
64) Phenanthrene-d10	17.151	188	1092132	20.000	ng	0.00
76) Chrysene-d12	21.386	240	1044209	20.000	ng	0.00
86) Perylene-d12	24.374	264	1027174	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.369	112	306899	19.895	ng	0.00
7) Phenol-d6	6.946	99	391479	19.279	ng	0.00
23) Nitrobenzene-d5	8.934	82	373577	18.841	ng	0.00
42) 2,4,6-Tribromophenol	15.898	330	127081	18.693	ng	0.00
45) 2-Fluorobiphenyl	13.039	172	918043	20.795	ng	0.00
79) Terphenyl-d14	19.780	244	1136477	20.625	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.293	88	71029	10.505	ng	97
3) Pyridine	3.687	79	173970	9.935	ng	98
4) n-Nitrosodimethylamine	3.599	42	34098	9.417	ng	# 99
6) Aniline	7.110	93	259330	9.660	ng	99
8) 2-Chlorophenol	7.351	128	161982	9.547	ng	99
9) Benzaldehyde	6.928	77	139011	10.767	ng	99
10) Phenol	6.969	94	211509	9.844	ng	99
11) bis(2-Chloroethyl)ether	7.210	93	169669	9.818	ng	99
12) 1,3-Dichlorobenzene	7.681	146	188081	9.915	ng	98
13) 1,4-Dichlorobenzene	7.822	146	198637	10.065	ng	99
14) 1,2-Dichlorobenzene	8.140	146	187370	9.959	ng	99
15) Benzyl Alcohol	8.022	79	134400	9.280	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.316	45	122149	10.175	ng	97
17) 2-Methylphenol	8.222	107	134124	9.461	ng	100
18) Hexachloroethane	8.869	117	70045	9.800	ng	98
19) n-Nitroso-di-n-propyla...	8.592	70	121867	9.807	ng	97
20) 3+4-Methylphenols	8.545	107	176847	9.324	ng	99
22) Acetophenone	8.604	105	260808	10.123	ng	# 98
24) Nitrobenzene	8.975	77	173945	9.646	ng	99
25) Isophorone	9.504	82	321944	9.407	ng	99
26) 2-Nitrophenol	9.687	139	64131	8.239	ng	96
27) 2,4-Dimethylphenol	9.751	122	153659	9.598	ng	99
28) bis(2-Chloroethoxy)met...	9.992	93	213076	9.531	ng	99
29) 2,4-Dichlorophenol	10.216	162	135013	9.122	ng	95
30) 1,2,4-Trichlorobenzene	10.439	180	158337	9.614	ng	99
31) Naphthalene	10.628	128	536516	10.063	ng	99
32) Benzoic acid	9.810	122	65412	6.508	ng	99
33) 4-Chloroaniline	10.728	127	218016	9.591	ng	99
34) Hexachlorobutadiene	10.922	225	94772	9.674	ng	98
35) Caprolactam	11.469	113	39590	8.437	ng	95
36) 4-Chloro-3-methylphenol	11.845	107	142482	9.022	ng	98
37) 2-Methylnaphthalene	12.239	142	305385	9.495	ng	97
38) 1-Methylnaphthalene	12.457	142	328309	9.617	ng	99
40) 1,2,4,5-Tetrachloroben...	12.610	216	167598	9.711	ng	98
41) Hexachlorocyclopentadiene	12.598	237	91203	8.701	ng	95
43) 2,4,6-Trichlorophenol	12.845	196	103556	9.124	ng	99

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44) 2,4,5-Trichlorophenol	12.910	196	114017	9.149	ng	98
46) 1,1'-Biphenyl	13.251	154	452133	10.102	ng	100
47) 2-Chloronaphthalene	13.286	162	356687	10.251	ng	99
48) 2-Nitroaniline	13.480	65	61491	8.031	ng	89
49) Acenaphthylene	14.133	152	548052	9.738	ng	99
50) Dimethylphthalate	13.874	163	384143	9.503	ng	99
51) 2,6-Dinitrotoluene	13.980	165	69112	8.478	ng	95
52) Acenaphthene	14.480	154	351485	9.985	ng	99
53) 3-Nitroaniline	14.304	138	75418	8.301	ng	97
54) 2,4-Dinitrophenol	14.510	184	27888	10.741	ng	98
55) Dibenzofuran	14.816	168	511101	9.925	ng	98
56) 4-Nitrophenol	14.604	139	64450	8.335	ng	99
57) 2,4-Dinitrotoluene	14.763	165	84481	7.856	ng	91
58) Fluorene	15.463	166	405716	10.286	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.033	232	98715m	9.156	ng	
60) Diethylphthalate	15.239	149	380099	9.477	ng	100
61) 4-Chlorophenyl-phenyle...	15.463	204	196034	10.283	ng	100
62) 4-Nitroaniline	15.463	138	70517	8.281	ng	97
63) Azobenzene	15.751	77	397398	9.913	ng	97
65) 4,6-Dinitro-2-methylph...	15.527	198	40699	6.823	ng	94
66) n-Nitrosodiphenylamine	15.668	169	341695	9.957	ng	99
67) 4-Bromophenyl-phenylether	16.351	248	111429	9.568	ng	96
68) Hexachlorobenzene	16.463	284	130925	9.622	ng	99
69) Atrazine	16.621	200	96720	8.846	ng	98
70) Pentachlorophenol	16.798	266	75315	8.513	ng	100
71) Phenanthrene	17.192	178	630983	10.114	ng	99
72) Anthracene	17.286	178	619530	9.926	ng	99
73) Carbazole	17.551	167	547511	9.629	ng	100
74) Di-n-butylphthalate	18.133	149	549702	8.885	ng	99
75) Fluoranthene	19.204	202	622189	9.457	ng	98
77) Benzidine	19.392	184	199082	6.780	ng	99
78) Pyrene	19.568	202	682132	9.692	ng	99
80) Butylbenzylphthalate	20.480	149	181298	7.723	ng	96
81) Benzo(a)anthracene	21.368	228	642785	9.725	ng	98
82) 3,3'-Dichlorobenzidine	21.292	252	157266	7.803	ng	99
83) Chrysene	21.427	228	618398	9.836	ng	99
84) Bis(2-ethylhexyl)phtha...	21.321	149	299821	8.291	ng	96
85) Di-n-octyl phthalate	22.462	149	380681	7.094	ng	99
87) Indeno(1,2,3-cd)pyrene	27.738	276	607722	9.189	ng	# 92
88) Benzo(b)fluoranthene	23.433	252	555747	9.306	ng	98
89) Benzo(k)fluoranthene	23.497	252	589937	9.677	ng	99
90) Benzo(a)pyrene	24.233	252	517175	9.211	ng	99
91) Dibenzo(a,h)anthracene	27.809	278	496324	9.245	ng	98
92) Benzo(g,h,i)perylene	28.785	276	502232	9.347	ng	98

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

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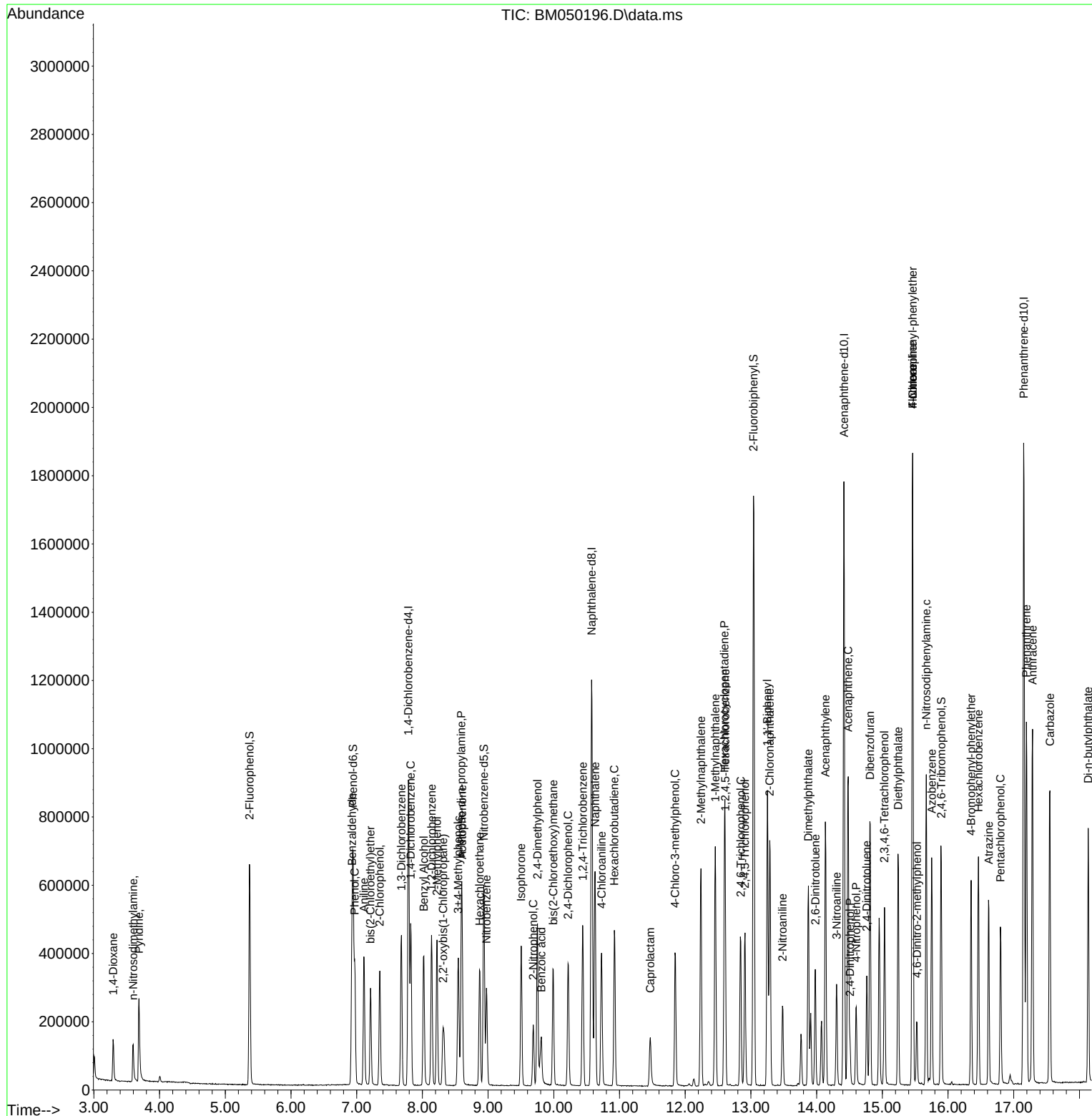
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