

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040825\
 Data File : BM049849.D
 Acq On : 08 Apr 2025 14:14
 Operator : RC/JU
 Sample : SSTDICC005
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC005

Manual Integrations
 APPROVED

Reviewed By :Anahy Claudio 04/09/2025
 Supervised By :Jagrut Upadhyay 04/09/2025

Quant Time: Apr 09 02:59:45 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM040825.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 09 02:53:19 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.781	152	315494	20.000	ng	0.00
21) Naphthalene-d8	10.575	136	1095995	20.000	ng	0.00
39) Acenaphthene-d10	14.421	164	671126	20.000	ng	0.00
64) Phenanthrene-d10	17.162	188	1305854	20.000	ng	0.00
76) Chrysene-d12	21.403	240	1255417	20.000	ng	0.00
86) Perylene-d12	24.403	264	1363674	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.363	112	180654	9.723	ng	0.00
7) Phenol-d6	6.951	99	219444	9.493	ng	0.00
23) Nitrobenzene-d5	8.933	82	183508	9.324	ng	0.00
42) 2,4,6-Tribromophenol	15.909	330	86574	8.869	ng	0.00
45) 2-Fluorobiphenyl	13.039	172	478149	9.661	ng	0.00
79) Terphenyl-d14	19.792	244	620191	9.258	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.263	88	39196	5.123	ng	99
3) Pyridine	3.669	79	99537	4.953	ng	93
4) n-Nitrosodimethylamine	3.575	42	37684	4.926	ng	# 91
6) Aniline	7.104	93	102863	5.193	ng	98
8) 2-Chlorophenol	7.345	128	93043	4.887	ng	99
9) Benzaldehyde	6.922	77	68896m	5.813	ng	
10) Phenol	6.975	94	108517	4.811	ng	100
11) bis(2-Chloroethyl)ether	7.204	93	87813	4.964	ng	99
12) 1,3-Dichlorobenzene	7.669	146	111770	4.965	ng	98
13) 1,4-Dichlorobenzene	7.816	146	111836	4.937	ng	99
14) 1,2-Dichlorobenzene	8.133	146	106307	4.983	ng	96
15) Benzyl Alcohol	8.022	79	67196	4.616	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.310	45	100016	5.073	ng	97
17) 2-Methylphenol	8.228	107	67662	4.867	ng	99
18) Hexachloroethane	8.863	117	39805	5.051	ng	96
19) n-Nitroso-di-n-propyla...	8.581	70	61968	4.840	ng	97
20) 3+4-Methylphenols	8.551	107	89903	4.744	ng	99
22) Acetophenone	8.598	105	126780	4.760	ng	96
24) Nitrobenzene	8.975	77	90100	4.754	ng	99
25) Isophorone	9.504	82	156018	4.732	ng	99
26) 2-Nitrophenol	9.686	139	44371	4.417	ng	96
27) 2,4-Dimethylphenol	9.757	122	50734	4.457	ng	98
28) bis(2-Chloroethoxy)met...	9.986	93	105841	4.795	ng	98
29) 2,4-Dichlorophenol	10.227	162	87444	4.621	ng	97
30) 1,2,4-Trichlorobenzene	10.439	180	105742	4.830	ng	99
31) Naphthalene	10.622	128	274967	4.883	ng	99
33) 4-Chloroaniline	10.733	127	95317	4.793	ng	99
34) Hexachlorobutadiene	10.916	225	65197	4.815	ng	98
35) Caprolactam	11.498	113	25608	4.718	ng	91
36) 4-Chloro-3-methylphenol	11.869	107	75972	4.583	ng	99
37) 2-Methylnaphthalene	12.239	142	187208	4.718	ng	99
38) 1-Methylnaphthalene	12.457	142	186332	4.820	ng	99
40) 1,2,4,5-Tetrachloroben...	12.610	216	111415	4.745	ng	99
41) Hexachlorocyclopentadiene	12.592	237	35937	4.368	ng	98
43) 2,4,6-Trichlorophenol	12.851	196	69935	4.681	ng	97
44) 2,4,5-Trichlorophenol	12.927	196	78011	4.805	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 1,1'-Biphenyl	13.257	154	243448	4.880	ng	99
47) 2-Chloronaphthalene	13.292	162	191299	4.879	ng	98
48) 2-Nitroaniline	13.498	65	38874	4.285	ng	94
49) Acenaphthylene	14.145	152	272990	4.779	ng	99
50) Dimethylphthalate	13.874	163	229200	4.841	ng	99
51) 2,6-Dinitrotoluene	13.992	165	41712	4.205	ng	93
52) Acenaphthene	14.486	154	172362	4.744	ng	100
53) 3-Nitroaniline	14.321	138	39627	4.133	ng	# 99
55) Dibenzofuran	14.821	168	291515	4.800	ng	99
56) 4-Nitrophenol	14.639	139	32685	3.826	ng	95
57) 2,4-Dinitrotoluene	14.780	165	51881	3.920	ng	# 94
58) Fluorene	15.468	166	227189	4.706	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.051	232	70065	4.924	ng	98
60) Diethylphthalate	15.239	149	220014	4.846	ng	98
61) 4-Chlorophenyl-phenyle...	15.468	204	117238	4.529	ng	97
62) 4-Nitroaniline	15.480	138	38105	3.843	ng	95
63) Azobenzene	15.757	77	180758	4.650	ng	98
66) n-Nitrosodiphenylamine	15.674	169	183131	4.686	ng	100
67) 4-Bromophenyl-phenylether	16.356	248	69182	4.564	ng	98
68) Hexachlorobenzene	16.474	284	81189	4.669	ng	95
69) Atrazine	16.627	200	60729	5.993	ng	98
70) Pentachlorophenol	16.821	266	59080	4.615	ng	98
71) Phenanthrene	17.203	178	336674	4.703	ng	100
72) Anthracene	17.298	178	325638	4.616	ng	99
73) Carbazole	17.568	167	301001	4.530	ng	99
74) Di-n-butylphthalate	18.133	149	365358	4.773	ng	100
75) Fluoranthene	19.221	202	387674	4.404	ng	100
77) Benzidine	19.409	184	100183	6.015	ng	97
78) Pyrene	19.586	202	411233	4.753	ng	99
80) Butylbenzylphthalate	20.486	149	157696	4.851	ng	98
81) Benzo(a)anthracene	21.386	228	388113	4.652	ng	100
82) 3,3'-Dichlorobenzidine	21.309	252	134300	4.262	ng	# 98
83) Chrysene	21.444	228	370038	4.665	ng	100
84) Bis(2-ethylhexyl)phtha...	21.315	149	237828	5.021	ng	100
85) Di-n-octyl phthalate	22.444	149	405083	4.889	ng	99
87) Indeno(1,2,3-cd)pyrene	27.785	276	468774	4.612	ng	# 93
88) Benzo(b)fluoranthene	23.456	252	389022	4.467	ng	98
89) Benzo(k)fluoranthene	23.521	252	385821	4.571	ng	98
90) Benzo(a)pyrene	24.262	252	344800	4.505	ng	98
91) Dibenzo(a,h)anthracene	27.844	278	384717	4.595	ng	99
92) Benzo(g,h,i)perylene	28.838	276	401539	4.693	ng	99

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

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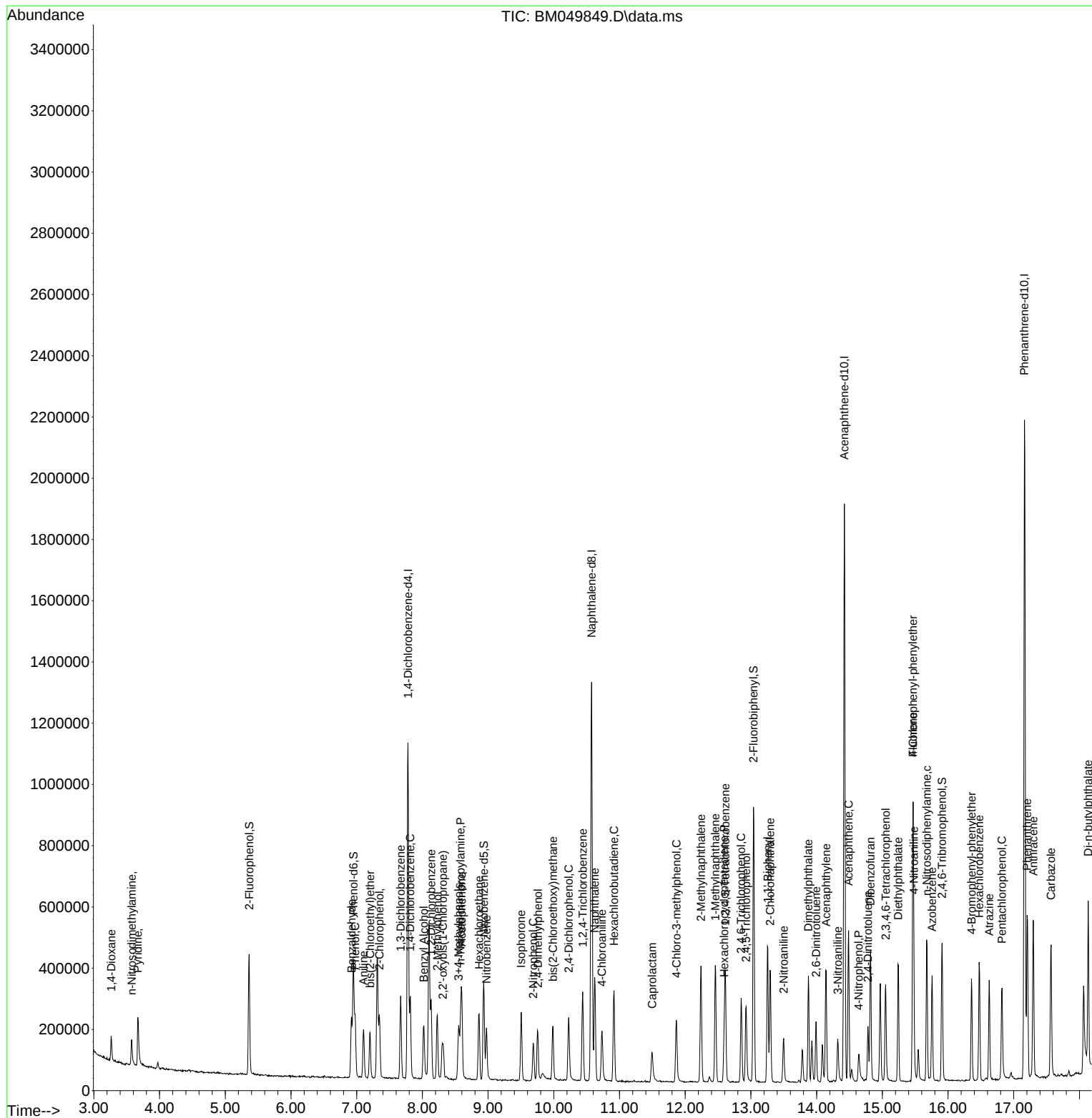
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