

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082024\
 Data File : BM047263.D
 Acq On : 20 Aug 2024 09:11
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC040

Quant Time: Aug 20 10:58:42 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	7.363	152	345222	20.000	ng	-0.01
21) Naphthalene-d8	10.116	136	1315377	20.000	ng	-0.02
39) Acenaphthene-d10	14.021	164	877836	20.000	ng	-0.01
64) Phenanthrene-d10	16.786	188	1861658	20.000	ng	0.00
76) Chrysene-d12	21.027	240	1404241	20.000	ng	0.00
86) Perylene-d12	23.727	264	1377858	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.016	112	1401926	72.563	ng	0.00
7) Phenol-d6	6.575	99	1896342	71.233	ng	-0.01
23) Nitrobenzene-d5	8.510	82	1865496	71.412	ng	-0.01
42) 2,4,6-Tribromophenol	15.533	330	846580	83.172	ng	0.00
45) 2-Fluorobiphenyl	12.633	172	4440513	78.028	ng	-0.01
79) Terphenyl-d14	19.456	244	6408452	97.788	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.028	88	278873	34.837	ng	96
3) Pyridine	3.405	79	777691	33.130	ng	97
4) n-Nitrosodimethylamine	3.322	42	332215	32.393	ng	97
6) Aniline	6.710	93	884405	35.148	ng	98
8) 2-Chlorophenol	6.940	128	775093	36.783	ng	97
9) Benzaldehyde	6.528	77	561185	40.308	ng	98
10) Phenol	6.604	94	940882	35.420	ng	98
11) bis(2-Chloroethyl)ether	6.816	93	795041	34.949	ng	98
12) 1,3-Dichlorobenzene	7.251	146	951345	37.993	ng	98
13) 1,4-Dichlorobenzene	7.398	146	956146	37.714	ng	99
14) 1,2-Dichlorobenzene	7.704	146	889745	37.195	ng	99
15) Benzyl Alcohol	7.616	79	679865	35.883	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.898	45	968400	33.175	ng	99
17) 2-Methylphenol	7.822	107	647365	36.529	ng	99
18) Hexachloroethane	8.410	117	361826	36.844	ng	94
19) n-Nitroso-di-n-propyla...	8.175	70	611695	34.112	ng	97
20) 3+4-Methylphenols	8.151	107	899910	36.323	ng	98
22) Acetophenone	8.181	105	1187745	37.408	ng	# 98
24) Nitrobenzene	8.551	77	962542	36.592	ng	98
25) Isophorone	9.081	82	1725309	37.932	ng	97
26) 2-Nitrophenol	9.251	139	490897	42.140	ng	96
27) 2,4-Dimethylphenol	9.334	122	537173	39.365	ng	99
28) bis(2-Chloroethoxy)met...	9.563	93	1077400	37.645	ng	99
29) 2,4-Dichlorophenol	9.786	162	854671	42.212	ng	99
30) 1,2,4-Trichlorobenzene	9.986	180	947706	41.360	ng	99
31) Naphthalene	10.169	128	2497949	38.121	ng	100
32) Benzoic acid	9.534	122	689176	43.471	ng	97
33) 4-Chloroaniline	10.292	127	953589	37.842	ng	100
34) Hexachlorobutadiene	10.451	225	597769	44.572	ng	100
35) Caprolactam	11.116	113	279113	39.695	ng	91
36) 4-Chloro-3-methylphenol	11.445	107	846919	39.276	ng	99
37) 2-Methylnaphthalene	11.792	142	1822700	39.067	ng	98
38) 1-Methylnaphthalene	12.022	142	1801652	39.073	ng	97
40) 1,2,4,5-Tetrachloroben...	12.175	216	1031986	43.101	ng	98
41) Hexachlorocyclopentadiene	12.151	237	317490	37.974	ng	96
43) 2,4,6-Trichlorophenol	12.439	196	717692	42.172	ng	99

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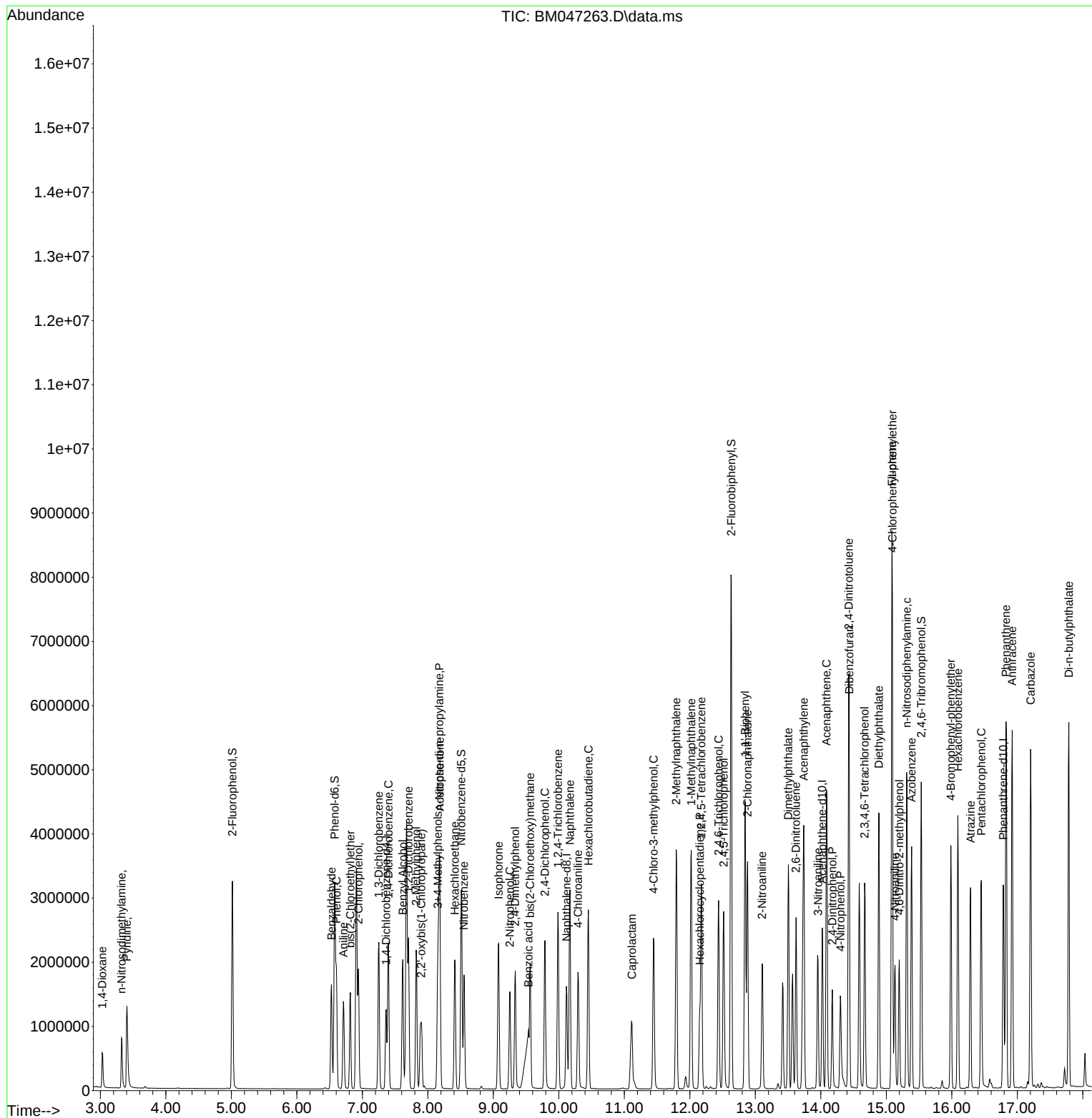
Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.516	196	792350	41.934	ng	97
46) 1,1'-Biphenyl	12.845	154	2386895	38.194	ng	99
47) 2-Chloronaphthalene	12.880	162	1850629	37.887	ng	99
48) 2-Nitroaniline	13.104	65	566000	36.987	ng	97
49) Acenaphthylene	13.739	152	2749552	38.188	ng	100
50) Dimethylphthalate	13.504	163	2399222	39.133	ng	99
51) 2,6-Dinitrotoluene	13.622	165	573841	40.040	ng	90
52) Acenaphthene	14.086	154	1733397	37.953	ng	100
53) 3-Nitroaniline	13.957	138	552457	38.469	ng	94
54) 2,4-Dinitrophenol	14.174	184	349854	40.730	ng	93
55) Dibenzofuran	14.433	168	2777357	38.499	ng	97
56) 4-Nitrophenol	14.298	139	458224	37.003	ng	98
57) 2,4-Dinitrotoluene	14.427	165	764192	39.949	ng	96
58) Fluorene	15.086	166	2269598	38.129	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.669	232	650127	41.865	ng	94
60) Diethylphthalate	14.886	149	2398331	38.308	ng	98
61) 4-Chlorophenyl-phenyle...	15.092	204	1142269	40.308	ng	94
62) 4-Nitroaniline	15.133	138	525113	37.116	ng	96
63) Azobenzene	15.386	77	2092385	35.248	ng	95
65) 4,6-Dinitro-2-methylph...	15.198	198	464873	42.650	ng	89
66) n-Nitrosodiphenylamine	15.310	169	1955591	38.567	ng	99
67) 4-Bromophenyl-phenylether	15.986	248	745098	41.535	ng	98
68) Hexachlorobenzene	16.092	284	867541	40.379	ng	96
69) Atrazine	16.286	200	584478	38.305	ng	98
70) Pentachlorophenol	16.451	266	618770	41.085	ng	99
71) Phenanthrene	16.827	178	3522062	37.698	ng	100
72) Anthracene	16.921	178	3467408	38.142	ng	100
73) Carbazole	17.204	167	3426865	36.920	ng	99
74) Di-n-butylphthalate	17.786	149	4190638	38.801	ng	99
75) Fluoranthene	18.868	202	4289860	37.749	ng	98
77) Benzidine	19.068	184	1172573	49.239	ng	99
78) Pyrene	19.233	202	4383390	43.717	ng	99
80) Butylbenzylphthalate	20.168	149	1854673	45.326	ng	96
81) Benzo(a)anthracene	21.015	228	3584675	38.453	ng	100
82) 3,3'-Dichlorobenzidine	20.950	252	1239869	42.509	ng	98
83) Chrysene	21.068	228	3340312	37.767	ng	99
84) Bis(2-ethylhexyl)phtha...	20.968	149	2554579	44.022	ng #	98
85) Di-n-octyl phthalate	21.997	149	4207980	42.663	ng	98
87) Indeno(1,2,3-cd)pyrene	26.727	276	3342517	37.530	ng	98
88) Benzo(b)fluoranthene	22.886	252	3168367	38.721	ng	99
89) Benzo(k)fluoranthene	22.939	252	3069657	39.717	ng	99
90) Benzo(a)pyrene	23.603	252	2735755	38.596	ng	99
91) Dibenzo(a,h)anthracene	26.774	278	2687520	37.288	ng	99
92) Benzo(g,h,i)perylene	27.662	276	2842276	37.982	ng	99

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

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