

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042925\
 Data File : BP024463.D
 Acq On : 29 Apr 2025 17:22
 Operator : RC/JU
 Sample : Q1871-04MS
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MH-AMS

Quant Time: Apr 29 18:03:30 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP041425.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 18 12:04:48 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.710	152	159274	20.000	ng	-0.01
21) Naphthalene-d8	10.481	136	640994	20.000	ng	#-0.02
39) Acenaphthene-d10	14.339	164	417040	20.000	ng	-0.01
64) Phenanthrene-d10	17.139	188	854293	20.000	ng	0.00
76) Chrysene-d12	21.598	240	948102	20.000	ng	# 0.00
86) Perylene-d12	24.944	264	1010012	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.340	112	1307321	135.716	ng	0.00
7) Phenol-d6	6.904	99	1620665	122.871	ng	0.00
23) Nitrobenzene-d5	8.857	82	1020350	90.768	ng	-0.01
42) 2,4,6-Tribromophenol	15.857	330	1015443	175.988	ng	0.00
45) 2-Fluorobiphenyl	12.951	172	2481491	91.105	ng	-0.01
79) Terphenyl-d14	19.874	244	4578147	97.330	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.293	88	148124	36.354	ng	# 64
3) Pyridine	3.687	79	335449	31.179	ng	94
4) n-Nitrosodimethylamine	3.587	42	167648	46.951	ng	77
6) Aniline	7.051	93	235189	19.954	ng	98
8) 2-Chlorophenol	7.287	128	506657	47.109	ng	99
9) Benzaldehyde	6.863	77	45857	7.028	ng	99
10) Phenol	6.928	94	586183	44.302	ng	91
11) bis(2-Chloroethyl)ether	7.145	93	428386	40.183	ng	99
12) 1,3-Dichlorobenzene	7.598	146	504772	43.596	ng	98
13) 1,4-Dichlorobenzene	7.745	146	515736	43.901	ng	98
14) 1,2-Dichlorobenzene	8.057	146	496999	43.812	ng	99
15) Benzyl Alcohol	7.951	79	428613	50.248	ng	97
16) 2,2'-oxybis(1-Chloropr...	8.228	45	477120	41.405	ng	100
17) 2-Methylphenol	8.157	107	422198	49.325	ng	97
18) Hexachloroethane	8.775	117	190469	44.864	ng	97
19) n-Nitroso-di-n-propyla...	8.510	70	338619	41.497	ng	94
20) 3+4-Methylphenols	8.481	107	575718	48.475	ng	98
22) Acetophenone	8.528	105	713343	44.963	ng	# 98
24) Nitrobenzene	8.898	77	505011	45.412	ng	99
25) Isophorone	9.422	82	993352	48.819	ng	97
26) 2-Nitrophenol	9.604	139	276980	50.896	ng	96
27) 2,4-Dimethylphenol	9.669	122	482496	70.610	ng	98
28) bis(2-Chloroethoxy)met...	9.898	93	588031	42.780	ng	99
29) 2,4-Dichlorophenol	10.145	162	490656	52.662	ng	99
30) 1,2,4-Trichlorobenzene	10.351	180	466993	46.781	ng	98
31) Naphthalene	10.533	128	1485996	44.010	ng	99
32) Benzoic acid	9.845	122	476683m	59.868	ng	
33) 4-Chloroaniline	10.651	127	82069	6.902	ng	97
34) Hexachlorobutadiene	10.816	225	296159	50.487	ng	100
35) Caprolactam	11.457	113	148999m	43.341	ng	
36) 4-Chloro-3-methylphenol	11.780	107	580849	53.523	ng	99
37) 2-Methylnaphthalene	12.139	142	967846	41.331	ng	97
38) 1-Methylnaphthalene	12.369	142	1025195	44.753	ng	98
40) 1,2,4,5-Tetrachloroben...	12.516	216	546119	47.618	ng	99
41) Hexachlorocyclopentadiene	12.492	237	749587	184.802	ng	99
43) 2,4,6-Trichlorophenol	12.763	196	410687	53.861	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.845	196	458737	54.329	ng	99
46) 1,1'-Biphenyl	13.163	154	1410427	46.010	ng	98
47) 2-Chloronaphthalene	13.204	162	1064510	46.463	ng	98
48) 2-Nitroaniline	13.416	65	346714	53.950	ng	96
49) Acenaphthylene	14.057	152	1835253	50.878	ng	100
50) Dimethylphthalate	13.798	163	1527155	50.355	ng	100
51) 2,6-Dinitrotoluene	13.916	165	334868	51.997	ng	95
52) Acenaphthene	14.404	154	1060509	46.375	ng	98
53) 3-Nitroaniline	14.251	138	126924	18.752	ng	97
54) 2,4-Dinitrophenol	14.463	184	408027	107.559	ng	90
55) Dibenzofuran	14.745	168	1704788	45.607	ng	97
56) 4-Nitrophenol	14.580	139	658821	112.701	ng	92
57) 2,4-Dinitrotoluene	14.716	165	488406	55.595	ng	98
58) Fluorene	15.404	166	1419218	49.046	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.980	232	425921	54.692	ng	99
60) Diethylphthalate	15.180	149	1581796	50.612	ng	100
61) 4-Chlorophenyl-phenyle...	15.404	204	703059	50.034	ng	99
62) 4-Nitroaniline	15.439	138	322268	44.976	ng	91
63) Azobenzene	15.698	77	1286506	44.786	ng	96
65) 4,6-Dinitro-2-methylph...	15.498	198	279414	51.878	ng	94
66) n-Nitrosodiphenylamine	15.621	169	1119056	43.887	ng	100
67) 4-Bromophenyl-phenylether	16.316	248	475695	50.913	ng	99
68) Hexachlorobenzene	16.433	284	588681	53.023	ng	99
69) Atrazine	16.604	200	224875	34.627	ng	99
70) Pentachlorophenol	16.786	266	851378	109.922	ng	99
71) Phenanthrene	17.180	178	2233026	47.915	ng	100
72) Anthracene	17.280	178	2239269	49.854	ng	99
73) Carbazole	17.562	167	2004400	45.671	ng	99
74) Di-n-butylphthalate	18.145	149	2862676	52.356	ng	99
75) Fluoranthene	19.274	202	2704191	48.788	ng	97
77) Benzidine	19.486	184	59652	4.964	ng	95
78) Pyrene	19.651	202	2782654	46.183	ng	99
80) Butylbenzylphthalate	20.604	149	1337546	51.827	ng	96
81) Benzo(a)anthracene	21.574	228	2864040	48.600	ng	100
82) 3,3'-Dichlorobenzidine	21.498	252	179894	8.697	ng	99
83) Chrysene	21.645	228	2680814	47.483	ng	100
84) Bis(2-ethylhexyl)phtha...	21.509	149	1974581	52.192	ng	99
85) Di-n-octyl phthalate	22.786	149	3334131	54.208	ng	100
87) Indeno(1,2,3-cd)pyrene	28.744	276	3981408m	56.780	ng	
88) Benzo(b)fluoranthene	23.886	252	3167956	52.327	ng	# 98
89) Benzo(k)fluoranthene	23.956	252	3018588	51.618	ng	# 97
90) Benzo(a)pyrene	24.792	252	2828702	54.166	ng	# 97
91) Dibenzo(a,h)anthracene	28.827	278	3284486	56.433	ng	97
92) Benzo(g,h,i)perylene	29.938	276	3152649	53.218	ng	96

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

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