

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052125\
 Data File : BP024739.D
 Acq On : 21 May 2025 16:09
 Operator : RC/JU
 Sample : PB168095BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB168095BS

Quant Time: May 21 16:30:47 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP051325.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 13 16:21:39 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.652	152	131263	20.000	ng	-0.01
21) Naphthalene-d8	10.416	136	507114	20.000	ng	-0.02
39) Acenaphthene-d10	14.281	164	323292	20.000	ng	-0.02
64) Phenanthrene-d10	17.081	188	610169	20.000	ng	-0.02
76) Chrysene-d12	21.516	240	717958	20.000	ng	0.00
86) Perylene-d12	24.780	264	839311	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.287	112	965182	125.764	ng	0.00
7) Phenol-d6	6.858	99	1137106	116.093	ng	0.00
23) Nitrobenzene-d5	8.805	82	699190	69.664	ng	0.00
42) 2,4,6-Tribromophenol	15.804	330	787700	143.722	ng	0.00
45) 2-Fluorobiphenyl	12.893	172	1874636	75.969	ng	-0.01
79) Terphenyl-d14	19.804	244	3387954	80.914	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.223	88	110320	30.613	ng	97
3) Pyridine	3.617	79	262732	32.879	ng	95
4) n-Nitrosodimethylamine	3.529	42	143737	40.745	ng	# 88
6) Aniline	6.993	93	295742	23.386	ng	99
8) 2-Chlorophenol	7.234	128	389118	44.663	ng	97
9) Benzaldehyde	6.811	77	209277	33.006	ng	98
10) Phenol	6.881	94	407641	40.323	ng	96
11) bis(2-Chloroethyl)ether	7.087	93	315913	38.030	ng	99
12) 1,3-Dichlorobenzene	7.546	146	410772	40.981	ng	98
13) 1,4-Dichlorobenzene	7.687	146	421630	41.823	ng	99
14) 1,2-Dichlorobenzene	7.999	146	399363	40.654	ng	99
15) Benzyl Alcohol	7.899	79	294271	39.535	ng	99
16) 2,2'-oxybis(1-Chloropr...	8.169	45	345225	34.702	ng	98
17) 2-Methylphenol	8.111	107	289936	39.987	ng	98
18) Hexachloroethane	8.717	117	157212	40.694	ng	98
19) n-Nitroso-di-n-propyla...	8.458	70	235766	34.797	ng	95
20) 3+4-Methylphenols	8.440	107	395800	40.129	ng	99
22) Acetophenone	8.469	105	527509	41.643	ng	99
24) Nitrobenzene	8.846	77	353263	39.996	ng	95
25) Isophorone	9.364	82	639758	36.747	ng	97
26) 2-Nitrophenol	9.546	139	208499	48.175	ng	97
27) 2,4-Dimethylphenol	9.622	122	337320	44.206	ng	98
28) bis(2-Chloroethoxy)met...	9.840	93	416471	40.040	ng	98
29) 2,4-Dichlorophenol	10.099	162	339875	45.963	ng	97
30) 1,2,4-Trichlorobenzene	10.287	180	372619	44.717	ng	97
31) Naphthalene	10.469	128	1105720	42.076	ng	100
32) Benzoic acid	9.805	122	215476	41.720	ng	94
33) 4-Chloroaniline	10.599	127	183013	17.258	ng	98
34) Hexachlorobutadiene	10.752	225	251681	46.624	ng	97
35) Caprolactam	11.393	113	111155	41.548	ng	98
36) 4-Chloro-3-methylphenol	11.734	107	394581	43.757	ng	99
37) 2-Methylnaphthalene	12.081	142	727486	42.753	ng	97
38) 1-Methylnaphthalene	12.305	142	761870	41.846	ng	100
40) 1,2,4,5-Tetrachloroben...	12.458	216	433678	46.195	ng	99
41) Hexachlorocyclopentadiene	12.434	237	597042	104.881	ng	97
43) 2,4,6-Trichlorophenol	12.710	196	294212	48.384	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.799	196	323501	47.794	ng	98
46) 1,1'-Biphenyl	13.105	154	1055344	44.166	ng	99
47) 2-Chloronaphthalene	13.146	162	800541	43.939	ng	99
48) 2-Nitroaniline	13.363	65	233432	43.276	ng	94
49) Acenaphthylene	13.999	152	1285084	42.149	ng	99
50) Dimethylphthalate	13.740	163	1036305	42.055	ng	99
51) 2,6-Dinitrotoluene	13.863	165	226632	43.548	ng	95
52) Acenaphthene	14.346	154	750344	42.809	ng	99
53) 3-Nitroaniline	14.210	138	123077	23.162	ng	95
54) 2,4-Dinitrophenol	14.416	184	269469	81.945	ng	98
55) Dibenzofuran	14.687	168	1204737	43.091	ng	98
56) 4-Nitrophenol	14.557	139	373188	81.061	ng	97
57) 2,4-Dinitrotoluene	14.663	165	339037	46.210	ng	98
58) Fluorene	15.346	166	988457	43.359	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.934	232	294256	46.864	ng	96
60) Diethylphthalate	15.122	149	1084714	43.553	ng	98
61) 4-Chlorophenyl-phenyle...	15.340	204	503437	44.282	ng	98
62) 4-Nitroaniline	15.387	138	223919m	43.642	ng	
63) Azobenzene	15.640	77	857948	40.268	ng	94
65) 4,6-Dinitro-2-methylph...	15.446	198	182840	46.391	ng	98
66) n-Nitrosodiphenylamine	15.563	169	863879	46.093	ng	99
67) 4-Bromophenyl-phenylether	16.251	248	354483	49.303	ng	96
68) Hexachlorobenzene	16.375	284	443741	48.570	ng	94
69) Atrazine	16.546	200	327372	47.663	ng	99
70) Pentachlorophenol	16.734	266	565169	110.201	ng	98
71) Phenanthrene	17.122	178	1491730	43.978	ng	100
72) Anthracene	17.216	178	1532702	44.911	ng	100
73) Carbazole	17.504	167	1395438	45.176	ng	100
74) Di-n-butylphthalate	18.087	149	1879200	47.305	ng	100
75) Fluoranthene	19.210	202	1856015	46.809	ng	98
77) Benzidine	19.410	184	773167	39.540	ng	99
78) Pyrene	19.587	202	1888163	43.892	ng	100
80) Butylbenzylphthalate	20.539	149	891758	46.611	ng	94
81) Benzo(a)anthracene	21.492	228	2014036	44.675	ng	99
82) 3,3'-Dichlorobenzidine	21.422	252	550613	32.884	ng	100
83) Chrysene	21.563	228	1886331	44.140	ng	100
84) Bis(2-ethylhexyl)phtha...	21.428	149	1312376	46.669	ng	99
85) Di-n-octyl phthalate	22.675	149	2208309	48.022	ng	99
87) Indeno(1,2,3-cd)pyrene	28.504	276	2788355	45.506	ng	# 88
88) Benzo(b)fluoranthene	23.757	252	2155576	44.870	ng	99
89) Benzo(k)fluoranthene	23.822	252	2228869	45.025	ng	99
90) Benzo(a)pyrene	24.633	252	2097566	45.462	ng	98
91) Dibenzo(a,h)anthracene	28.563	278	2306937	46.279	ng	98
92) Benzo(g,h,i)perylene	29.651	276	2222348	44.831	ng	97

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

