

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052225\
 Data File : BP024756.D
 Acq On : 22 May 2025 12:58
 Operator : RC/JU
 Sample : PB168098BSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB168098BSD

Quant Time: May 22 13:25:32 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	126327	20.000	ng	-0.01
21) Naphthalene-d8	10.416	136	523024	20.000	ng	-0.02
39) Acenaphthene-d10	14.287	164	338733	20.000	ng	-0.02
64) Phenanthrene-d10	17.092	188	686645	20.000	ng	0.00
76) Chrysene-d12	21.510	240	768545	20.000	ng	-0.01
86) Perylene-d12	24.786	264	835982	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.287	112	1006990	136.339	ng	0.00
7) Phenol-d6	6.858	99	1245584	132.137	ng	0.00
23) Nitrobenzene-d5	8.805	82	764802	73.883	ng	0.00
42) 2,4,6-Tribromophenol	15.810	330	848083	147.685	ng	0.00
45) 2-Fluorobiphenyl	12.899	172	1985294	76.786	ng	0.00
79) Terphenyl-d14	19.816	244	3652739	81.496	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.223	88	107291	30.935	ng	98
3) Pyridine	3.617	79	272244	35.401	ng	97
4) n-Nitrosodimethylamine	3.528	42	142518	41.978	ng	95
6) Aniline	6.993	93	278550	22.887	ng	99
8) 2-Chlorophenol	7.234	128	400340	47.747	ng	99
9) Benzaldehyde	6.811	77	221040	36.224	ng	97
10) Phenol	6.881	94	447784	46.024	ng	98
11) bis(2-Chloroethyl)ether	7.087	93	321348	40.196	ng	99
12) 1,3-Dichlorobenzene	7.546	146	411266	42.633	ng	98
13) 1,4-Dichlorobenzene	7.687	146	421327	43.426	ng	98
14) 1,2-Dichlorobenzene	7.999	146	408503	43.209	ng	99
15) Benzyl Alcohol	7.899	79	325668	45.463	ng	98
16) 2,2'-oxybis(1-Chloropr...	8.169	45	379444	39.632	ng	99
17) 2-Methylphenol	8.111	107	324116	46.447	ng	98
18) Hexachloroethane	8.716	117	154783	41.631	ng	98
19) n-Nitroso-di-n-propyla...	8.452	70	254858	39.085	ng	96
20) 3+4-Methylphenols	8.440	107	431650	45.474	ng	98
22) Acetophenone	8.469	105	553115	42.336	ng	# 99
24) Nitrobenzene	8.846	77	382330	41.971	ng	94
25) Isophorone	9.369	82	732635	40.802	ng	97
26) 2-Nitrophenol	9.546	139	218042	48.847	ng	98
27) 2,4-Dimethylphenol	9.622	122	369301	46.925	ng	100
28) bis(2-Chloroethoxy)met...	9.840	93	441817	41.185	ng	99
29) 2,4-Dichlorophenol	10.093	162	375212	49.198	ng	97
30) 1,2,4-Trichlorobenzene	10.287	180	383813	44.659	ng	98
31) Naphthalene	10.469	128	1176453	43.406	ng	100
32) Benzoic acid	9.810	122	270871m	49.215	ng	
33) 4-Chloroaniline	10.605	127	165916	15.170	ng	99
34) Hexachlorobutadiene	10.752	225	250650	45.021	ng	96
35) Caprolactam	11.387	113	126254	45.756	ng	98
36) 4-Chloro-3-methylphenol	11.740	107	436668	46.951	ng	99
37) 2-Methylnaphthalene	12.081	142	753865	42.956	ng	99
38) 1-Methylnaphthalene	12.304	142	807085	42.981	ng	100
40) 1,2,4,5-Tetrachloroben...	12.463	216	446029	45.345	ng	100
41) Hexachlorocyclopentadiene	12.434	237	507943	85.162	ng	97
43) 2,4,6-Trichlorophenol	12.710	196	314461	49.357	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.804	196	349258	49.247	ng	96
46) 1,1'-Biphenyl	13.104	154	1090713	43.566	ng	99
47) 2-Chloronaphthalene	13.152	162	833303	43.653	ng	99
48) 2-Nitroaniline	13.363	65	249151	44.085	ng	96
49) Acenaphthylene	14.004	152	1391550	43.560	ng	99
50) Dimethylphthalate	13.746	163	1142952	44.269	ng	100
51) 2,6-Dinitrotoluene	13.869	165	247920	45.467	ng	99
52) Acenaphthene	14.351	154	799545	43.536	ng	99
53) 3-Nitroaniline	14.210	138	117478	21.100	ng	96
54) 2,4-Dinitrophenol	14.422	184	322278	92.631	ng	96
55) Dibenzofuran	14.693	168	1288076	43.972	ng	100
56) 4-Nitrophenol	14.563	139	402969	83.411	ng	97
57) 2,4-Dinitrotoluene	14.669	165	365228	47.511	ng	97
58) Fluorene	15.351	166	1071652	44.866	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.934	232	317896	48.321	ng	98
60) Diethylphthalate	15.128	149	1184185	45.380	ng	99
61) 4-Chlorophenyl-phenyle...	15.345	204	547042	45.924	ng	100
62) 4-Nitroaniline	15.393	138	254227m	47.290	ng	
63) Azobenzene	15.645	77	949950	42.554	ng	95
65) 4,6-Dinitro-2-methylph...	15.451	198	218741	49.318	ng	97
66) n-Nitrosodiphenylamine	15.569	169	961369	45.582	ng	99
67) 4-Bromophenyl-phenylether	16.257	248	378264	46.751	ng	97
68) Hexachlorobenzene	16.387	284	480414	46.728	ng	94
69) Atrazine	16.551	200	373961	48.382	ng	98
70) Pentachlorophenol	16.745	266	621802	107.740	ng	99
71) Phenanthrene	17.134	178	1679605	44.002	ng	100
72) Anthracene	17.228	178	1732101	45.101	ng	99
73) Carbazole	17.516	167	1619014	46.577	ng	100
74) Di-n-butylphthalate	18.104	149	2111662	47.236	ng	100
75) Fluoranthene	19.222	202	2044320	45.816	ng	99
77) Benzidine	19.422	184	742781	35.485	ng	99
78) Pyrene	19.592	202	2099503	45.592	ng	99
80) Butylbenzylphthalate	20.545	149	992487	48.461	ng	95
81) Benzo(a)anthracene	21.492	228	2194524	45.475	ng	99
82) 3,3'-Dichlorobenzidine	21.428	252	544789	30.394	ng	99
83) Chrysene	21.557	228	2031513	44.408	ng	100
84) Bis(2-ethylhexyl)phtha...	21.428	149	1478644	49.121	ng	99
85) Di-n-octyl phthalate	22.675	149	2442323	49.615	ng	99
87) Indeno(1,2,3-cd)pyrene	28.515	276	2735441	44.820	ng	# 92
88) Benzo(b)fluoranthene	23.763	252	2233590	46.679	ng	99
89) Benzo(k)fluoranthene	23.827	252	2248794	45.608	ng	99
90) Benzo(a)pyrene	24.639	252	2120304	46.138	ng	98
91) Dibenzo(a,h)anthracene	28.598	278	2246534	45.246	ng	99
92) Benzo(g,h,i)perylene	29.668	276	2167806	43.905	ng	97

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

