

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111225\
 Data File : BF144236.D
 Acq On : 12 Nov 2025 18:50
 Operator : RC/JU
 Sample : PB170515BS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB170515BS

Manual Integrations
 APPROVED

Reviewed By :Jagrut
 Upadhyay

Quant Time: Nov 13 10:41:37 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 05 18:37:33 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.869	152	45402	20.000	ng	0.00
21) Naphthalene-d8	8.151	136	180251	20.000	ng	# 0.00
39) Acenaphthene-d10	9.910	164	101053	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	166539	20.000	ng	0.00
76) Chrysene-d12	14.045	240	84699	20.000	ng	# 0.00
86) Perylene-d12	15.527	264	112671	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	334570	115.316	ng	0.01
7) Phenol-d6	6.504	99	385516	117.270	ng	0.00
23) Nitrobenzene-d5	7.434	82	249688	80.004	ng	0.00
42) 2,4,6-Tribromophenol	10.704	330	139884	110.310	ng	0.00
45) 2-Fluorobiphenyl	9.228	172	509312	74.438	ng	0.00
79) Terphenyl-d14	12.986	244	500617	93.883	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.728	88	50178	41.833	ng	98
3) Pyridine	3.487	79	147170	43.978	ng	99
4) n-Nitrosodimethylamine	3.440	42	84384	48.275	ng	92
6) Aniline	6.534	93	153296	32.729	ng	96
8) 2-Chlorophenol	6.657	128	124556	43.776	ng	98
9) Benzaldehyde	6.422	77	75921	40.898	ng	98
10) Phenol	6.516	94	167739	41.762	ng	93
11) bis(2-Chloroethyl)ether	6.604	93	134789	46.055	ng	97
12) 1,3-Dichlorobenzene	6.810	146	150921	42.985	ng	96
13) 1,4-Dichlorobenzene	6.886	146	149097	42.388	ng	99
14) 1,2-Dichlorobenzene	7.039	146	141537	41.981	ng	99
15) Benzyl Alcohol	7.010	79	116903	44.909	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.139	45	268794	49.873	ng	98
17) 2-Methylphenol	7.122	107	101915	45.025	ng	97
18) Hexachloroethane	7.381	117	49609	42.451	ng	99
19) n-Nitroso-di-n-propyla...	7.281	70	94930	43.865	ng	97
20) 3+4-Methylphenols	7.275	107	109917	41.194	ng	97
22) Acetophenone	7.275	105	157378	40.700	ng	93
24) Nitrobenzene	7.451	77	144126	42.532	ng	99
25) Isophorone	7.686	82	263493	44.635	ng	100
26) 2-Nitrophenol	7.769	139	74133	44.194	ng	98
27) 2,4-Dimethylphenol	7.804	122	119044	47.349	ng	99
28) bis(2-Chloroethoxy)met...	7.898	93	161394	43.781	ng	97
29) 2,4-Dichlorophenol	8.010	162	117076	40.402	ng	99
30) 1,2,4-Trichlorobenzene	8.092	180	123719	41.578	ng	98
31) Naphthalene	8.175	128	346477	40.409	ng	99
32) Benzoic acid	7.933	122	79816	43.309	ng	94
33) 4-Chloroaniline	8.222	127	51152	16.357	ng	97
34) Hexachlorobutadiene	8.286	225	85612	38.176	ng	99
35) Caprolactam	8.598	113	38426m	48.401	ng	
36) 4-Chloro-3-methylphenol	8.710	107	112760	43.419	ng	98
37) 2-Methylnaphthalene	8.863	142	214401	37.399	ng	99
38) 1-Methylnaphthalene	8.963	142	219790	39.734	ng	100
40) 1,2,4,5-Tetrachloroben...	9.033	216	125809	37.988	ng	98
41) Hexachlorocyclopentadiene	9.016	237	115457	86.467	ng	98
43) 2,4,6-Trichlorophenol	9.145	196	93067	41.288	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.186	196	92917	38.247	ng	97
46) 1,1'-Biphenyl	9.328	154	285865	40.526	ng	99
47) 2-Chloronaphthalene	9.357	162	240609	39.988	ng	100
48) 2-Nitroaniline	9.451	65	77854	44.844	ng	100
49) Acenaphthylene	9.775	152	370278	45.159	ng	100
50) Dimethylphthalate	9.633	163	285446	42.638	ng	99
51) 2,6-Dinitrotoluene	9.698	165	62667	46.242	ng	96
52) Acenaphthene	9.945	154	206698	37.697	ng	98
53) 3-Nitroaniline	9.863	138	37825	25.533	ng	98
54) 2,4-Dinitrophenol	9.975	184	56433	68.960	ng	97
55) Dibenzofuran	10.116	168	306185	39.871	ng	99
56) 4-Nitrophenol	10.033	139	85624	79.902	ng	98
57) 2,4-Dinitrotoluene	10.104	165	79721	43.943	ng	91
58) Fluorene	10.463	166	235506	40.336	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.239	232	73860	42.971	ng	98
60) Diethylphthalate	10.333	149	265856	43.079	ng	99
61) 4-Chlorophenyl-phenyle...	10.451	204	131377	39.502	ng	99
62) 4-Nitroaniline	10.480	138	59767	42.364	ng	93
63) Azobenzene	10.610	77	263433	45.866	ng	99
65) 4,6-Dinitro-2-methylph...	10.516	198	45609	40.313	ng	97
66) n-Nitrosodiphenylamine	10.569	169	241641	45.113	ng	99
67) 4-Bromophenyl-phenylether	10.939	248	86208	40.555	ng	96
68) Hexachlorobenzene	11.010	284	103591	41.373	ng	98
69) Atrazine	11.098	200	79812	46.854	ng	98
70) Pentachlorophenol	11.204	266	97885	78.509	ng	97
71) Phenanthrene	11.427	178	351466	42.388	ng	100
72) Anthracene	11.474	178	358443	42.513	ng	100
73) Carbazole	11.633	167	305936	42.672	ng	99
74) Di-n-butylphthalate	11.957	149	399956	46.147	ng	99
75) Fluoranthene	12.616	202	344156	40.180	ng	98
77) Benzidine	12.739	184	155632	62.120	ng	99
78) Pyrene	12.845	202	332878	48.743	ng	100
80) Butylbenzylphthalate	13.463	149	124612	52.647	ng	100
81) Benzo(a)anthracene	14.033	228	273075	46.518	ng	99
82) 3,3'-Dichlorobenzidine	13.998	252	50489	25.092	ng	94
83) Chrysene	14.074	228	241482	44.562	ng	99
84) Bis(2-ethylhexyl)phtha...	14.015	149	158832	49.020	ng	100
85) Di-n-octyl phthalate	14.633	149	241505	43.200	ng	98
87) Indeno(1,2,3-cd)pyrene	17.027	276	334616	41.372	ng	97
88) Benzo(b)fluoranthene	15.092	252	298592m	46.614	ng	
89) Benzo(k)fluoranthene	15.121	252	249896	41.680	ng	98
90) Benzo(a)pyrene	15.462	252	270772	45.821	ng	# 97
91) Dibenzo(a,h)anthracene	17.045	278	270183	41.595	ng	99
92) Benzo(g,h,i)perylene	17.480	276	266444	41.538	ng	97

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

