

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111925\
 Data File : BF144281.D
 Acq On : 19 Nov 2025 16:51
 Operator : RC/JU
 Sample : PB170637BS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB170637BS

Manual Integrations
 APPROVED

Reviewed By :Jagrut
 Upadhyay

Quant Time: Nov 19 17:10:28 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	56226	20.000	ng	0.00
21) Naphthalene-d8	8.151	136	216046	20.000	ng	0.00
39) Acenaphthene-d10	9.910	164	113709	20.000	ng	0.00
64) Phenanthrene-d10	11.398	188	177131	20.000	ng	0.00
76) Chrysene-d12	14.051	240	137709	20.000	ng	0.00
86) Perylene-d12	15.527	264	163785	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.498	112	429166	119.445	ng	0.00
7) Phenol-d6	6.504	99	471025	115.698	ng	0.00
23) Nitrobenzene-d5	7.434	82	311791	83.350	ng	0.00
42) 2,4,6-Tribromophenol	10.704	330	160761	112.663	ng	0.00
45) 2-Fluorobiphenyl	9.228	172	650048	84.433	ng	0.00
79) Terphenyl-d14	12.986	244	614080	70.831	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.687	88	60746	40.894	ng	99
3) Pyridine	3.457	79	170483	41.137	ng	99
4) n-Nitrosodimethylamine	3.404	42	100647	46.494	ng	95
6) Aniline	6.534	93	179106	30.878	ng	94
8) 2-Chlorophenol	6.657	128	159537	45.277	ng	97
9) Benzaldehyde	6.422	77	79878	34.746	ng	99
10) Phenol	6.522	94	212487	42.719	ng	82
11) bis(2-Chloroethyl)ether	6.604	93	170533	47.051	ng	99
12) 1,3-Dichlorobenzene	6.810	146	199727	45.935	ng	97
13) 1,4-Dichlorobenzene	6.887	146	201249	46.200	ng	100
14) 1,2-Dichlorobenzene	7.039	146	192617	46.134	ng	98
15) Benzyl Alcohol	7.010	79	151103	46.873	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.139	45	308352	46.199	ng	92
17) 2-Methylphenol	7.122	107	125951	44.932	ng	97
18) Hexachloroethane	7.381	117	65664	45.373	ng	97
19) n-Nitroso-di-n-propyla...	7.281	70	118923	44.373	ng	98
20) 3+4-Methylphenols	7.275	107	145185	43.937	ng	# 89
22) Acetophenone	7.275	105	232218	50.105	ng	97
24) Nitrobenzene	7.451	77	182598	44.958	ng	99
25) Isophorone	7.686	82	324337	45.838	ng	100
26) 2-Nitrophenol	7.769	139	94217	46.861	ng	96
27) 2,4-Dimethylphenol	7.804	122	143819	47.725	ng	98
28) bis(2-Chloroethoxy)met...	7.898	93	202859	45.912	ng	98
29) 2,4-Dichlorophenol	8.010	162	152425	43.886	ng	100
30) 1,2,4-Trichlorobenzene	8.092	180	162139	45.462	ng	99
31) Naphthalene	8.175	128	450116	43.798	ng	99
32) Benzoic acid	7.928	122	101455	45.930	ng	93
33) 4-Chloroaniline	8.222	127	65954	17.596	ng	99
34) Hexachlorobutadiene	8.286	225	114067	42.438	ng	99
35) Caprolactam	8.592	113	44691m	46.966	ng	
36) 4-Chloro-3-methylphenol	8.710	107	136965	44.001	ng	100
37) 2-Methylnaphthalene	8.863	142	274770	39.989	ng	100
38) 1-Methylnaphthalene	8.963	142	273635	41.273	ng	98
40) 1,2,4,5-Tetrachloroben...	9.033	216	186297	49.991	ng	99
41) Hexachlorocyclopentadiene	9.016	237	106161	74.867	ng	99
43) 2,4,6-Trichlorophenol	9.145	196	116268	45.840	ng	100

11/20/2025
 Supervised By :Sohil
 Odhani

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44) 2,4,5-Trichlorophenol	9.186	196	122461	44.797	ng	96
46) 1,1'-Biphenyl	9.328	154	396489	49.953	ng	100
47) 2-Chloronaphthalene	9.357	162	301744	44.567	ng	100
48) 2-Nitroaniline	9.451	65	92724	47.465	ng	97
49) Acenaphthylene	9.769	152	466180	50.527	ng	100
50) Dimethylphthalate	9.628	163	360391	47.842	ng	100
51) 2,6-Dinitrotoluene	9.692	165	76134	49.927	ng	99
52) Acenaphthene	9.945	154	253419	41.074	ng	99
53) 3-Nitroaniline	9.863	138	51834	31.095	ng	98
54) 2,4-Dinitrophenol	9.975	184	50294	54.618	ng	100
55) Dibenzofuran	10.116	168	376764	43.601	ng	100
56) 4-Nitrophenol	10.033	139	89842	74.506	ng	97
57) 2,4-Dinitrotoluene	10.098	165	99158	48.573	ng	98
58) Fluorene	10.457	166	294797	44.871	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.233	232	87051	45.009	ng	98
60) Diethylphthalate	10.328	149	328012	47.235	ng	98
61) 4-Chlorophenyl-phenyle...	10.451	204	167429	44.739	ng	99
62) 4-Nitroaniline	10.480	138	66395	41.824	ng	94
63) Azobenzene	10.610	77	295429	45.712	ng	98
65) 4,6-Dinitro-2-methylph...	10.510	198	44038	36.597	ng	98
66) n-Nitrosodiphenylamine	10.569	169	283857	49.826	ng	99
67) 4-Bromophenyl-phenylether	10.939	248	107051	47.349	ng	97
68) Hexachlorobenzene	11.010	284	122468	45.988	ng	96
69) Atrazine	11.098	200	93234	51.460	ng	99
70) Pentachlorophenol	11.204	266	108952	82.160	ng	99
71) Phenanthrene	11.422	178	400296	45.390	ng	100
72) Anthracene	11.474	178	410935	45.825	ng	99
73) Carbazole	11.633	167	353141	46.311	ng	99
74) Di-n-butylphthalate	11.957	149	439794	47.709	ng	99
75) Fluoranthene	12.616	202	410843	45.097	ng	99
77) Benzidine	12.739	184	156609	38.447	ng	99
78) Pyrene	12.845	202	420865	37.904	ng	99
80) Butylbenzylphthalate	13.457	149	166102	43.162	ng	99
81) Benzo(a)anthracene	14.039	228	453626	47.529	ng	99
82) 3,3'-Dichlorobenzidine	13.998	252	116378	35.573	ng	96
83) Chrysene	14.074	228	417293	47.363	ng	99
84) Bis(2-ethylhexyl)phtha...	14.015	149	232499	44.134	ng	100
85) Di-n-octyl phthalate	14.633	149	424090	46.659	ng	100
87) Indeno(1,2,3-cd)pyrene	17.033	276	419473	35.678	ng	98
88) Benzo(b)fluoranthene	15.098	252	479083	51.451	ng	99
89) Benzo(k)fluoranthene	15.127	252	445276	51.090	ng	99
90) Benzo(a)pyrene	15.468	252	447750	52.123	ng	100
91) Dibenzo(a,h)anthracene	17.045	278	342769	36.302	ng	98
92) Benzo(g,h,i)perylene	17.486	276	311583	33.416	ng	98

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

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Abundance TIC: BF144281.D\data.ms

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