

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080125\
 Data File : BF143278.D
 Acq On : 01 Aug 2025 13:03
 Operator : RC/JU
 Sample : PB169077BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB169077BS

Quant Time: Aug 01 13:25:27 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF071725.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 17 15:14:05 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.963	152	102726	20.000	ng	0.00
21) Naphthalene-d8	8.245	136	402894	20.000	ng	0.00
39) Acenaphthene-d10	9.998	164	212350	20.000	ng	0.00
64) Phenanthrene-d10	11.486	188	364957	20.000	ng	0.00
76) Chrysene-d12	14.127	240	204179	20.000	ng	0.00
86) Perylene-d12	15.633	264	207302	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.604	112	745968	114.914	ng	0.02
7) Phenol-d6	6.593	99	929722	113.786	ng	0.00
23) Nitrobenzene-d5	7.522	82	638897	79.046	ng	0.00
42) 2,4,6-Tribromophenol	10.792	330	291491	132.877	ng	0.00
45) 2-Fluorobiphenyl	9.316	172	1190626	76.681	ng	0.00
79) Terphenyl-d14	13.069	244	1164765	84.891	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.899	88	116179	37.823	ng	100
3) Pyridine	3.652	79	282017	35.373	ng	98
4) n-Nitrosodimethylamine	3.593	42	167772	40.758	ng	98
6) Aniline	6.622	93	373032	32.758	ng	94
8) 2-Chlorophenol	6.751	128	292350	42.965	ng	97
9) Benzaldehyde	6.510	77	184796	30.161	ng	98
10) Phenol	6.604	94	376841	42.336	ng	90
11) bis(2-Chloroethyl)ether	6.693	93	279114	40.954	ng	98
12) 1,3-Dichlorobenzene	6.904	146	308643	41.755	ng	99
13) 1,4-Dichlorobenzene	6.981	146	316003	42.451	ng	100
14) 1,2-Dichlorobenzene	7.134	146	301217	42.544	ng	99
15) Benzyl Alcohol	7.098	79	270636	43.379	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.234	45	509535	40.302	ng	98
17) 2-Methylphenol	7.210	107	246221	43.036	ng	98
18) Hexachloroethane	7.475	117	112825	43.782	ng	98
19) n-Nitroso-di-n-propyla...	7.369	70	214772	40.970	ng	99
20) 3+4-Methylphenols	7.363	107	295847	42.620	ng	# 76
22) Acetophenone	7.369	105	395536	41.467	ng	# 98
24) Nitrobenzene	7.540	77	322680	42.821	ng	99
25) Isophorone	7.781	82	591945	41.486	ng	99
26) 2-Nitrophenol	7.857	139	158096	48.341	ng	99
27) 2,4-Dimethylphenol	7.893	122	272687	40.905	ng	100
28) bis(2-Chloroethoxy)met...	7.987	93	353461	41.317	ng	100
29) 2,4-Dichlorophenol	8.098	162	244553	43.500	ng	100
30) 1,2,4-Trichlorobenzene	8.181	180	262022	43.141	ng	99
31) Naphthalene	8.263	128	839310	42.300	ng	100
32) Benzoic acid	8.004	122	153676	41.687	ng	100
33) 4-Chloroaniline	8.304	127	172008	21.230	ng	99
34) Hexachlorobutadiene	8.387	225	162144	43.702	ng	99
35) Caprolactam	8.675	113	82355m	48.477	ng	
36) 4-Chloro-3-methylphenol	8.787	107	270471	44.263	ng	100
37) 2-Methylnaphthalene	8.957	142	512864	42.476	ng	100
38) 1-Methylnaphthalene	9.057	142	530392	42.659	ng	99
40) 1,2,4,5-Tetrachloroben...	9.128	216	260264	42.452	ng	99
41) Hexachlorocyclopentadiene	9.116	237	328567	82.366	ng	99
43) 2,4,6-Trichlorophenol	9.234	196	176732	41.652	ng	99

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44) 2,4,5-Trichlorophenol	9.275	196	190710	44.933	ng	99
46) 1,1'-Biphenyl	9.422	154	711802	42.964	ng	99
47) 2-Chloronaphthalene	9.445	162	523882	42.736	ng	99
48) 2-Nitroaniline	9.534	65	181505	47.214	ng	100
49) Acenaphthylene	9.863	152	886064	43.131	ng	100
50) Dimethylphthalate	9.716	163	626516	45.263	ng	100
51) 2,6-Dinitrotoluene	9.775	165	137755	48.922	ng	99
52) Acenaphthene	10.034	154	566786	46.679	ng	98
53) 3-Nitroaniline	9.945	138	101661	30.867	ng	98
54) 2,4-Dinitrophenol	10.051	184	129077	99.382	ng	# 1
55) Dibenzofuran	10.204	168	770251	42.726	ng	99
56) 4-Nitrophenol	10.104	139	216601	88.070	ng	98
57) 2,4-Dinitrotoluene	10.181	165	185348	52.467	ng	95
58) Fluorene	10.551	166	597126	44.133	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.322	232	159905	45.482	ng	98
60) Diethylphthalate	10.416	149	627943	45.899	ng	100
61) 4-Chlorophenyl-phenyle...	10.539	204	299207	44.412	ng	100
62) 4-Nitroaniline	10.563	138	138303	48.996	ng	98
63) Azobenzene	10.698	77	615705	43.181	ng	99
65) 4,6-Dinitro-2-methylph...	10.592	198	91050	47.636	ng	98
66) n-Nitrosodiphenylamine	10.657	169	535107	41.638	ng	99
67) 4-Bromophenyl-phenylether	11.028	248	187731	43.163	ng	98
68) Hexachlorobenzene	11.104	284	196304	43.182	ng	96
69) Atrazine	11.181	200	171872	48.511	ng	99
70) Pentachlorophenol	11.292	266	225132	84.212	ng	98
71) Phenanthrene	11.516	178	834843	43.082	ng	100
72) Anthracene	11.563	178	845596	42.786	ng	100
73) Carbazole	11.716	167	763818	44.260	ng	100
74) Di-n-butylphthalate	12.039	149	921317	47.182	ng	100
75) Fluoranthene	12.698	202	833975	46.888	ng	99
77) Benzidine	12.816	184	337780	42.939	ng	98
78) Pyrene	12.933	202	816421	46.852	ng	99
80) Butylbenzylphthalate	13.539	149	299874	56.199	ng	99
81) Benzo(a)anthracene	14.116	228	633000	46.306	ng	99
82) 3,3'-Dichlorobenzidine	14.074	252	122772	26.947	ng	99
83) Chrysene	14.157	228	533282	43.390	ng	99
84) Bis(2-ethylhexyl)phtha...	14.098	149	387790	48.015	ng	100
85) Di-n-octyl phthalate	14.716	149	620118	42.359	ng	99
87) Indeno(1,2,3-cd)pyrene	17.180	276	660034	45.153	ng	97
88) Benzo(b)fluoranthene	15.186	252	565613	44.662	ng	100
89) Benzo(k)fluoranthene	15.221	252	510964	45.214	ng	98
90) Benzo(a)pyrene	15.568	252	529530	45.382	ng	100
91) Dibenzo(a,h)anthracene	17.198	278	535886	45.041	ng	98
92) Benzo(g,h,i)perylene	17.645	276	521256	45.291	ng	98

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

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