

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080125\
 Data File : BF143281.D
 Acq On : 01 Aug 2025 14:37
 Operator : RC/JU
 Sample : PB169087BS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB169087BS

Quant Time: Aug 01 15:08:44 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	111937	20.000	ng	0.00
21) Naphthalene-d8	8.245	136	478481	20.000	ng	0.00
39) Acenaphthene-d10	9.998	164	196679	20.000	ng	0.00
64) Phenanthrene-d10	11.486	188	332732	20.000	ng	0.00
76) Chrysene-d12	14.127	240	185103	20.000	ng	0.00
86) Perylene-d12	15.633	264	196335	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.604	112	846345	119.649	ng	0.02
7) Phenol-d6	6.592	99	1084828	121.844	ng	0.00
23) Nitrobenzene-d5	7.522	82	750639	78.200	ng	0.00
42) 2,4,6-Tribromophenol	10.792	330	279755	137.688	ng	0.00
45) 2-Fluorobiphenyl	9.316	172	1190719	82.798	ng	0.00
79) Terphenyl-d14	13.068	244	1102522	88.636	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.899	88	137544	41.094	ng	98
3) Pyridine	3.651	79	397516	45.758	ng	97
4) n-Nitrosodimethylamine	3.593	42	205667	45.853	ng	98
6) Aniline	6.622	93	362043	29.177	ng	95
8) 2-Chlorophenol	6.751	128	365340	49.273	ng	97
9) Benzaldehyde	6.510	77	212086	31.767	ng	99
10) Phenol	6.604	94	426617	43.984	ng	94
11) bis(2-Chloroethyl)ether	6.692	93	326661	43.986	ng	99
12) 1,3-Dichlorobenzene	6.904	146	354530	44.016	ng	99
13) 1,4-Dichlorobenzene	6.981	146	353562	43.588	ng	100
14) 1,2-Dichlorobenzene	7.134	146	344809	44.694	ng	99
15) Benzyl Alcohol	7.098	79	321266	47.257	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.234	45	627582	45.555	ng	98
17) 2-Methylphenol	7.210	107	284459	45.628	ng	99
18) Hexachloroethane	7.475	117	129864	46.247	ng	97
19) n-Nitroso-di-n-propyla...	7.369	70	243183	42.573	ng	100
20) 3+4-Methylphenols	7.363	107	334201	44.184	ng	# 74
22) Acetophenone	7.363	105	444791	39.264	ng	95
24) Nitrobenzene	7.539	77	372746	41.651	ng	99
25) Isophorone	7.775	82	686428	40.508	ng	99
26) 2-Nitrophenol	7.857	139	182404	46.963	ng	96
27) 2,4-Dimethylphenol	7.892	122	328255	41.462	ng	99
28) bis(2-Chloroethoxy)met...	7.986	93	463479	45.619	ng	99
29) 2,4-Dichlorophenol	8.098	162	306253	45.869	ng	99
30) 1,2,4-Trichlorobenzene	8.181	180	325305	45.099	ng	99
31) Naphthalene	8.263	128	1058902	44.936	ng	100
32) Benzoic acid	8.004	122	199637	45.114	ng	100
33) 4-Chloroaniline	8.304	127	122027	12.682	ng	99
34) Hexachlorobutadiene	8.386	225	195150	44.289	ng	98
35) Caprolactam	8.675	113	101664m	50.389	ng	
36) 4-Chloro-3-methylphenol	8.792	107	337599	46.521	ng	98
37) 2-Methylnaphthalene	8.957	142	648907	45.253	ng	99
38) 1-Methylnaphthalene	9.057	142	668891	45.299	ng	100
40) 1,2,4,5-Tetrachloroben...	9.128	216	308822	54.386	ng	99
41) Hexachlorocyclopentadiene	9.116	237	389339	105.377	ng	100
43) 2,4,6-Trichlorophenol	9.233	196	194434	49.476	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.275	196	209348	53.255	ng	99
46) 1,1'-Biphenyl	9.422	154	695805	45.345	ng	99
47) 2-Chloronaphthalene	9.445	162	510087	44.926	ng	99
48) 2-Nitroaniline	9.533	65	172299	48.390	ng	100
49) Acenaphthylene	9.863	152	864483	45.433	ng	100
50) Dimethylphthalate	9.716	163	604078	47.120	ng	100
51) 2,6-Dinitrotoluene	9.775	165	130014	49.852	ng	99
52) Acenaphthene	10.033	154	555470	49.392	ng	99
53) 3-Nitroaniline	9.945	138	73630	24.137	ng	98
54) 2,4-Dinitrophenol	10.051	184	121594	100.966	ng	# 1
55) Dibenzofuran	10.204	168	747478	44.767	ng	99
56) 4-Nitrophenol	10.104	139	199667	87.654	ng	99
57) 2,4-Dinitrotoluene	10.180	165	175695	53.697	ng	95
58) Fluorene	10.551	166	570237	45.504	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.322	232	153666	47.190	ng	97
60) Diethylphthalate	10.416	149	597644	47.165	ng	100
61) 4-Chlorophenyl-phenyle...	10.539	204	288333	46.208	ng	99
62) 4-Nitroaniline	10.557	138	127551	48.788	ng	98
63) Azobenzene	10.698	77	586851	44.437	ng	99
65) 4,6-Dinitro-2-methylph...	10.592	198	85778	49.044	ng	97
66) n-Nitrosodiphenylamine	10.657	169	510935	43.608	ng	99
67) 4-Bromophenyl-phenylether	11.027	248	177500	44.764	ng	99
68) Hexachlorobenzene	11.104	284	187073	45.137	ng	98
69) Atrazine	11.180	200	157449	48.744	ng	100
70) Pentachlorophenol	11.292	266	208299	85.462	ng	99
71) Phenanthrene	11.516	178	788552	44.634	ng	100
72) Anthracene	11.563	178	808098	44.848	ng	99
73) Carbazole	11.716	167	712207	45.267	ng	100
74) Di-n-butylphthalate	12.039	149	856407	48.106	ng	100
75) Fluoranthene	12.698	202	777745	47.962	ng	100
77) Benzidine	12.816	184	226771	31.798	ng	99
78) Pyrene	12.933	202	759881	48.101	ng	99
80) Butylbenzylphthalate	13.539	149	281240	58.138	ng	100
81) Benzo(a)anthracene	14.115	228	565267	45.613	ng	100
82) 3,3'-Dichlorobenzidine	14.074	252	88646	21.462	ng	98
83) Chrysene	14.151	228	522696	46.912	ng	99
84) Bis(2-ethylhexyl)phtha...	14.098	149	363243	49.610	ng	100
85) Di-n-octyl phthalate	14.715	149	586041	44.156	ng	98
87) Indeno(1,2,3-cd)pyrene	17.180	276	646594	46.705	ng	97
88) Benzo(b)fluoranthene	15.186	252	552673	46.078	ng	99
89) Benzo(k)fluoranthene	15.221	252	492088	45.976	ng	98
90) Benzo(a)pyrene	15.568	252	511924	46.324	ng	99
91) Dibenzo(a,h)anthracene	17.198	278	527401	46.804	ng	97
92) Benzo(g,h,i)perylene	17.645	276	607779	55.758	ng	98

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

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