

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031125\
 Data File : BF141910.D
 Acq On : 11 Mar 2025 11:51
 Operator : RC/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Mar 11 12:36:25 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF031025.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 10 15:46:22 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.881	152	212930	20.000	ng	0.00
21) Naphthalene-d8	8.163	136	840081	20.000	ng	0.00
39) Acenaphthene-d10	9.916	164	479107	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	819898	20.000	ng	0.00
76) Chrysene-d12	14.039	240	519479	20.000	ng	0.00
86) Perylene-d12	15.509	264	442953	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.492	112	966380	75.745	ng	0.00
7) Phenol-d6	6.504	99	1228433	75.624	ng	0.00
23) Nitrobenzene-d5	7.445	82	1166556	78.146	ng	0.00
42) 2,4,6-Tribromophenol	10.704	330	488429	80.359	ng	0.00
45) 2-Fluorobiphenyl	9.239	172	2374971	75.388	ng	0.00
79) Terphenyl-d14	12.986	244	2775945	79.004	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.675	88	207900	38.891	ng	99
3) Pyridine	3.434	79	509090	38.824	ng	99
4) n-Nitrosodimethylamine	3.393	42	242195	38.747	ng	98
6) Aniline	6.545	93	619097	38.634	ng	100
8) 2-Chlorophenol	6.669	128	542074	38.309	ng	99
9) Benzaldehyde	6.434	77	337776	37.180	ng	100
10) Phenol	6.516	94	651072	38.098	ng	98
11) bis(2-Chloroethyl)ether	6.616	93	484388	37.748	ng	100
12) 1,3-Dichlorobenzene	6.822	146	582982	38.163	ng	100
13) 1,4-Dichlorobenzene	6.898	146	594118	38.438	ng	100
14) 1,2-Dichlorobenzene	7.051	146	554766	38.106	ng	99
15) Benzyl Alcohol	7.022	79	506215	38.547	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.157	45	568739	36.712	ng	99
17) 2-Methylphenol	7.128	107	430396	38.255	ng	99
18) Hexachloroethane	7.392	117	225190	38.853	ng	100
19) n-Nitroso-di-n-propyla...	7.298	70	387430	37.118	ng	100
20) 3+4-Methylphenols	7.281	107	548388	38.055	ng	98
22) Acetophenone	7.292	105	763464	37.949	ng	99
24) Nitrobenzene	7.463	77	576817	38.869	ng	100
25) Isophorone	7.704	82	1009454	38.255	ng	100
26) 2-Nitrophenol	7.775	139	298908	41.039	ng	98
27) 2,4-Dimethylphenol	7.810	122	392258	39.112	ng	99
28) bis(2-Chloroethoxy)met...	7.910	93	629638	38.044	ng	100
29) 2,4-Dichlorophenol	8.016	162	481689	38.695	ng	100
30) 1,2,4-Trichlorobenzene	8.104	180	529203	38.575	ng	98
31) Naphthalene	8.186	128	1643492	38.085	ng	100
32) Benzoic acid	7.928	122	369497	41.912	ng	98
33) 4-Chloroaniline	8.233	127	585849	38.457	ng	99
34) Hexachlorobutadiene	8.304	225	351479	39.286	ng	98
35) Caprolactam	8.604	113	144643	38.112	ng	96
36) 4-Chloro-3-methylphenol	8.704	107	534161	38.467	ng	99
37) 2-Methylnaphthalene	8.875	142	1115480	38.672	ng	100
38) 1-Methylnaphthalene	8.975	142	1060457	38.099	ng	99
40) 1,2,4,5-Tetrachloroben...	9.039	216	568573	38.978	ng	98
41) Hexachlorocyclopentadiene	9.028	237	230480	40.373	ng	100
43) 2,4,6-Trichlorophenol	9.145	196	382004	40.071	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.186	196	372824	38.614	ng	100
46) 1,1'-Biphenyl	9.339	154	1402423	38.525	ng	100
47) 2-Chloronaphthalene	9.363	162	1049133	38.666	ng	99
48) 2-Nitroaniline	9.457	65	307247	39.109	ng	99
49) Acenaphthylene	9.780	152	1554840	38.616	ng	100
50) Dimethylphthalate	9.639	163	1283810	38.265	ng	100
51) 2,6-Dinitrotoluene	9.698	165	279974	40.079	ng	97
52) Acenaphthene	9.951	154	1091559	38.576	ng	99
53) 3-Nitroaniline	9.869	138	277307	39.135	ng	98
54) 2,4-Dinitrophenol	9.969	184	136926	38.237	ng	91
55) Dibenzofuran	10.122	168	1548206	38.292	ng	100
56) 4-Nitrophenol	10.016	139	212052	39.683	ng	99
57) 2,4-Dinitrotoluene	10.104	165	364725	39.975	ng	100
58) Fluorene	10.469	166	1187102	38.073	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.239	232	351000	39.950	ng	98
60) Diethylphthalate	10.339	149	1281977	38.320	ng	100
61) 4-Chlorophenyl-phenyle...	10.457	204	629512	38.726	ng	98
62) 4-Nitroaniline	10.480	138	268529	38.925	ng	99
63) Azobenzene	10.616	77	1194591	38.408	ng	100
65) 4,6-Dinitro-2-methylph...	10.510	198	195453	39.988	ng	99
66) n-Nitrosodiphenylamine	10.574	169	1014883	38.251	ng	100
67) 4-Bromophenyl-phenylether	10.945	248	401588	39.001	ng	100
68) Hexachlorobenzene	11.010	284	448381	39.703	ng	99
69) Atrazine	11.098	200	299767	36.855	ng	100
70) Pentachlorophenol	11.204	266	282858	40.651	ng	99
71) Phenanthrene	11.427	178	1684816	38.045	ng	100
72) Anthracene	11.480	178	1690241	38.043	ng	100
73) Carbazole	11.633	167	1449478	37.829	ng	100
74) Di-n-butylphthalate	11.963	149	1929199	37.945	ng	100
75) Fluoranthene	12.616	202	1750151	37.256	ng	100
77) Benzidine	12.733	184	392482	54.153	ng	99
78) Pyrene	12.845	202	1726338	38.418	ng	100
80) Butylbenzylphthalate	13.463	149	695571	39.548	ng	98
81) Benzo(a)anthracene	14.027	228	1266188	37.144	ng	100
82) 3,3'-Dichlorobenzidine	13.992	252	380363	38.611	ng	98
83) Chrysene	14.068	228	1214561	39.306	ng	99
84) Bis(2-ethylhexyl)phtha...	14.015	149	956586	39.447	ng	99
85) Di-n-octyl phthalate	14.633	149	1358481	40.262	ng	99
87) Indeno(1,2,3-cd)pyrene	16.998	276	1098821	38.348	ng	99
88) Benzo(b)fluoranthene	15.080	252	1084408	35.721	ng	100
89) Benzo(k)fluoranthene	15.109	252	1011237	38.924	ng	99
90) Benzo(a)pyrene	15.445	252	903590	38.039	ng	99
91) Dibenzo(a,h)anthracene	17.015	278	904778	38.270	ng	99
92) Benzo(g,h,i)perylene	17.445	276	885987	37.923	ng	99

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

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