

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF031025\
 Data File : BF141905.D
 Acq On : 10 Mar 2025 15:20
 Operator : RC/JU
 Sample : SSTDICC050
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC050

Quant Time: Mar 10 15:39:09 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 15:01:52 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.881	152	215017	20.000	ng	0.00
21) Naphthalene-d8	8.163	136	877329	20.000	ng	0.00
39) Acenaphthene-d10	9.922	164	481971	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	811142	20.000	ng	0.00
76) Chrysene-d12	14.045	240	491943	20.000	ng	0.00
86) Perylene-d12	15.509	264	413564	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.492	112	1249601	96.967	ng	0.00
7) Phenol-d6	6.504	99	1594608	97.156	ng	0.00
23) Nitrobenzene-d5	7.445	82	1507471	96.304	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	625035	102.342	ng	0.00
45) 2-Fluorobiphenyl	9.239	172	3064989	96.836	ng	0.00
79) Terphenyl-d14	12.992	244	3394803	102.007	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.675	88	266045	49.258	ng	99
3) Pyridine	3.434	79	646627	48.716	ng	99
4) n-Nitrosodimethylamine	3.398	42	320008	50.718	ng	99
6) Aniline	6.545	93	775854	47.884	ng	99
8) 2-Chlorophenol	6.669	128	696154	48.738	ng	100
9) Benzaldehyde	6.434	77	418049	45.569	ng	99
10) Phenol	6.522	94	829086	47.876	ng	96
11) bis(2-Chloroethyl)ether	6.616	93	631516	48.672	ng	100
12) 1,3-Dichlorobenzene	6.822	146	753352	48.807	ng	99
13) 1,4-Dichlorobenzene	6.898	146	756875	48.465	ng	100
14) 1,2-Dichlorobenzene	7.051	146	706635	48.049	ng	100
15) Benzyl Alcohol	7.022	79	652454	49.113	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.157	45	745710	47.687	ng	100
17) 2-Methylphenol	7.128	107	557981	49.153	ng	99
18) Hexachloroethane	7.392	117	290127	49.551	ng	99
19) n-Nitroso-di-n-propyla...	7.298	70	505166	47.887	ng	99
20) 3+4-Methylphenols	7.286	107	704563	48.388	ng	# 86
22) Acetophenone	7.292	105	987577	46.924	ng	97
24) Nitrobenzene	7.463	77	752936	48.514	ng	99
25) Isophorone	7.704	82	1327254	48.027	ng	99
26) 2-Nitrophenol	7.781	139	392772	51.644	ng	100
27) 2,4-Dimethylphenol	7.810	122	510395	48.725	ng	99
28) bis(2-Chloroethoxy)met...	7.910	93	826776	47.757	ng	100
29) 2,4-Dichlorophenol	8.016	162	647319	49.793	ng	100
30) 1,2,4-Trichlorobenzene	8.104	180	697274	48.617	ng	99
31) Naphthalene	8.186	128	2196111	48.863	ng	100
32) Benzoic acid	7.939	122	500593	54.676	ng	100
33) 4-Chloroaniline	8.233	127	775699	48.909	ng	99
34) Hexachlorobutadiene	8.304	225	469787	50.376	ng	99
35) Caprolactam	8.616	113	193267	48.891	ng	95
36) 4-Chloro-3-methylphenol	8.710	107	721800	49.925	ng	100
37) 2-Methylnaphthalene	8.875	142	1436485	47.638	ng	100
38) 1-Methylnaphthalene	8.975	142	1357356	46.482	ng	100
40) 1,2,4,5-Tetrachloroben...	9.039	216	719431	48.955	ng	99
41) Hexachlorocyclopentadiene	9.027	237	284968	49.142	ng	99
43) 2,4,6-Trichlorophenol	9.151	196	481782	50.291	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.186	196	482570	49.580	ng	99
46) 1,1'-Biphenyl	9.339	154	1809746	49.526	ng	99
47) 2-Chloronaphthalene	9.369	162	1360528	49.960	ng	99
48) 2-Nitroaniline	9.457	65	429111	54.854	ng	97
49) Acenaphthylene	9.780	152	2005538	49.674	ng	100
50) Dimethylphthalate	9.645	163	1716995	51.273	ng	100
51) 2,6-Dinitrotoluene	9.704	165	370368	53.184	ng	96
52) Acenaphthene	9.957	154	1377901	48.333	ng	98
53) 3-Nitroaniline	9.874	138	340633	47.638	ng	100
54) 2,4-Dinitrophenol	9.975	184	184410	49.104	ng	# 58
55) Dibenzofuran	10.127	168	1936780	47.647	ng	99
56) 4-Nitrophenol	10.022	139	276487	51.474	ng	99
57) 2,4-Dinitrotoluene	10.104	165	461276	50.321	ng	97
58) Fluorene	10.469	166	1497123	47.708	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.239	232	438025	49.600	ng	100
60) Diethylphthalate	10.339	149	1635073	48.639	ng	100
61) 4-Chlorophenyl-phenyle...	10.457	204	799744	48.929	ng	99
62) 4-Nitroaniline	10.486	138	346378	49.990	ng	99
63) Azobenzene	10.622	77	1519356	48.653	ng	98
65) 4,6-Dinitro-2-methylph...	10.516	198	262910	54.673	ng	99
66) n-Nitrosodiphenylamine	10.580	169	1293745	49.373	ng	100
67) 4-Bromophenyl-phenylether	10.951	248	532762	52.740	ng	97
68) Hexachlorobenzene	11.016	284	573467	51.359	ng	97
69) Atrazine	11.104	200	418147	65.441	ng	99
70) Pentachlorophenol	11.204	266	365292	53.050	ng	99
71) Phenanthrene	11.433	178	2080633	47.434	ng	100
72) Anthracene	11.480	178	2081490	47.299	ng	99
73) Carbazole	11.633	167	1831424	48.338	ng	99
74) Di-n-butylphthalate	11.963	149	2510447	50.185	ng	99
75) Fluoranthene	12.616	202	2328357	50.564	ng	99
77) Benzidine	12.733	184	270893	38.966	ng	99
78) Pyrene	12.845	202	2308671	54.723	ng	100
80) Butylbenzylphthalate	13.463	149	864849	52.022	ng	99
81) Benzo(a)anthracene	14.033	228	1592035	49.192	ng	100
82) 3,3'-Dichlorobenzidine	13.992	252	449408	48.044	ng	99
83) Chrysene	14.068	228	1465389	50.404	ng	100
84) Bis(2-ethylhexyl)phtha...	14.021	149	1203021	52.747	ng	99
85) Di-n-octyl phthalate	14.633	149	1692206	53.307	ng	100
87) Indeno(1,2,3-cd)pyrene	17.003	276	1366974	50.969	ng	99
88) Benzo(b)fluoranthene	15.080	252	1392083	49.625	ng	99
89) Benzo(k)fluoranthene	15.109	252	1167351	47.638	ng	100
90) Benzo(a)pyrene	15.451	252	1119256	50.591	ng	99
91) Dibenzo(a,h)anthracene	17.021	278	1131875	51.149	ng	99
92) Benzo(g,h,i)perylene	17.456	276	1099760	50.106	ng	99

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

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