

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF051625\
 Data File : BF142428.D
 Acq On : 16 May 2025 15:46
 Operator : RC/JU
 Sample : Q2034-01MS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 L3-WC-1MS

Quant Time: May 16 16:18:07 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF050525.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 05 18:41:44 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.904	152	137288	20.000	ng	0.00
21) Naphthalene-d8	8.186	136	518623	20.000	ng	0.00
39) Acenaphthene-d10	9.939	164	275083	20.000	ng	0.00
64) Phenanthrene-d10	11.427	188	447181	20.000	ng	0.00
76) Chrysene-d12	14.068	240	265603	20.000	ng	0.00
86) Perylene-d12	15.557	264	306381	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	569985	70.874	ng	0.01
7) Phenol-d6	6.522	99	807010	81.587	ng	0.00
23) Nitrobenzene-d5	7.463	82	400666	47.116	ng	0.00
42) 2,4,6-Tribromophenol	10.727	330	247945	93.357	ng	0.00
45) 2-Fluorobiphenyl	9.257	172	947431	49.110	ng	0.00
79) Terphenyl-d14	13.010	244	769053	42.856	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.746	88	136436	37.673	ng	99
3) Pyridine	3.499	79	360854	42.439	ng	97
4) n-Nitrosodimethylamine	3.451	42	197836	40.033	ng	97
6) Aniline	6.563	93	381647	39.337	ng	99
8) 2-Chlorophenol	6.687	128	383031	44.569	ng	97
9) Benzaldehyde	6.451	77	240611	41.562	ng	99
10) Phenol	6.540	94	473413	45.151	ng	98
11) bis(2-Chloroethyl)ether	6.634	93	344481	33.316	ng	98
12) 1,3-Dichlorobenzene	6.845	146	411895	41.893	ng	98
13) 1,4-Dichlorobenzene	6.922	146	414557	41.994	ng	99
14) 1,2-Dichlorobenzene	7.075	146	402024	42.427	ng	99
15) Benzyl Alcohol	7.040	79	304949	47.611	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.175	45	589098	38.153	ng	99
17) 2-Methylphenol	7.151	107	299914	43.105	ng	99
18) Hexachloroethane	7.416	117	144224	43.910	ng	99
19) n-Nitroso-di-n-propyla...	7.310	70	246907	40.355	ng	99
20) 3+4-Methylphenols	7.304	107	369929	42.601	ng	# 82
22) Acetophenone	7.310	105	487894	42.310	ng	99
24) Nitrobenzene	7.481	77	371972	46.404	ng	97
25) Isophorone	7.722	82	674365	41.878	ng	99
26) 2-Nitrophenol	7.798	139	200917	51.660	ng	99
27) 2,4-Dimethylphenol	7.834	122	355524	43.912	ng	99
28) bis(2-Chloroethoxy)met...	7.934	93	420387	40.599	ng	99
29) 2,4-Dichlorophenol	8.039	162	319814	46.463	ng	99
30) 1,2,4-Trichlorobenzene	8.128	180	335910	43.363	ng	99
31) Naphthalene	8.210	128	1087792	43.277	ng	99
32) Benzoic acid	7.934	122	215135	49.175	ng	99
33) 4-Chloroaniline	8.251	127	105510	10.287	ng	98
34) Hexachlorobutadiene	8.322	225	203670	44.401	ng	99
35) Caprolactam	8.616	113	102506	51.098	ng	97
36) 4-Chloro-3-methylphenol	8.733	107	322665	45.095	ng	99
37) 2-Methylnaphthalene	8.898	142	665741	42.625	ng	98
38) 1-Methylnaphthalene	8.998	142	691715	42.492	ng	99
40) 1,2,4,5-Tetrachloroben...	9.063	216	335611	43.683	ng	99
41) Hexachlorocyclopentadiene	9.051	237	427169	89.965	ng	100
43) 2,4,6-Trichlorophenol	9.175	196	238487	49.213	ng	100

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.210	196	241004	46.740	ng	99
46) 1,1'-Biphenyl	9.363	154	908286	43.064	ng	99
47) 2-Chloronaphthalene	9.386	162	676168	43.108	ng	100
48) 2-Nitroaniline	9.481	65	208366	51.826	ng	93
49) Acenaphthylene	9.804	152	1142094	43.526	ng	100
50) Dimethylphthalate	9.657	163	788021	44.612	ng	100
51) 2,6-Dinitrotoluene	9.722	165	175826	53.178	ng	89
52) Acenaphthene	9.975	154	759039	48.109	ng	100
53) 3-Nitroaniline	9.886	138	91461	23.586	ng	95
54) 2,4-Dinitrophenol	9.992	184	184155	103.927	ng	# 4
55) Dibenzofuran	10.145	168	1000094	42.951	ng	100
56) 4-Nitrophenol	10.045	139	285607	97.862	ng	95
57) 2,4-Dinitrotoluene	10.122	165	236805	55.835	ng	95
58) Fluorene	10.486	166	768569	43.251	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.263	232	203047	47.702	ng	98
60) Diethylphthalate	10.357	149	760296	43.100	ng	100
61) 4-Chlorophenyl-phenyle...	10.480	204	369512	43.101	ng	99
62) 4-Nitroaniline	10.504	138	173027	56.299	ng	99
63) Azobenzene	10.639	77	762104	45.211	ng	97
65) 4,6-Dinitro-2-methylph...	10.528	198	115734	54.395	ng	97
66) n-Nitrosodiphenylamine	10.598	169	675610	45.402	ng	100
67) 4-Bromophenyl-phenylether	10.969	248	230116	45.405	ng	98
68) Hexachlorobenzene	11.033	284	253071	45.396	ng	99
69) Atrazine	11.122	200	192964	49.194	ng	99
70) Pentachlorophenol	11.227	266	276241	94.193	ng	98
71) Phenanthrene	11.451	178	1030186	43.950	ng	100
72) Anthracene	11.504	178	1032598	42.874	ng	100
73) Carbazole	11.657	167	870883	40.304	ng	99
74) Di-n-butylphthalate	11.986	149	993331	43.718	ng	99
75) Fluoranthene	12.639	202	912620	38.949	ng	99
77) Benzidine	12.757	184	361702	74.937	ng	98
78) Pyrene	12.869	202	899478	38.047	ng	99
80) Butylbenzylphthalate	13.486	149	312036	45.438	ng	98
81) Benzo(a)anthracene	14.057	228	774173	45.008	ng	100
82) 3,3'-Dichlorobenzidine	14.016	252	149949	37.487	ng	100
83) Chrysene	14.092	228	710454	44.368	ng	99
84) Bis(2-ethylhexyl)phtha...	14.045	149	403147	45.146	ng	100
85) Di-n-octyl phthalate	14.657	149	705396	44.272	ng	97
87) Indeno(1,2,3-cd)pyrene	17.062	276	645400	30.282	ng	99
88) Benzo(b)fluoranthene	15.121	252	825541	44.489	ng	98
89) Benzo(k)fluoranthene	15.151	252	760594	44.790	ng	100
90) Benzo(a)pyrene	15.492	252	766334	46.086	ng	100
91) Dibenzo(a,h)anthracene	17.080	278	516680	29.344	ng	99
92) Benzo(g,h,i)perylene	17.515	276	475436	27.206	ng	97

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

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