

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082824\
 Data File : BF139202.D
 Acq On : 28 Aug 2024 11:38
 Operator : RC/JU
 Sample : PB163034BSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB163034BSD

Quant Time: Aug 28 12:29:17 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 27 02:51:47 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.893	152	105164	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	395427	20.000	ng	0.00
39) Acenaphthene-d10	9.928	164	218164	20.000	ng	0.00
64) Phenanthrene-d10	11.416	188	395038	20.000	ng	0.00
76) Chrysene-d12	14.051	240	193809	20.000	ng	0.00
86) Perylene-d12	15.521	264	187033	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.522	112	612233	97.825	ng	0.02
7) Phenol-d6	6.522	99	800166	93.160	ng	0.00
23) Nitrobenzene-d5	7.457	82	785861	110.101	ng	0.00
42) 2,4,6-Tribromophenol	10.716	330	219863	117.670	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	1311147	93.355	ng	0.00
79) Terphenyl-d14	12.998	244	1347470	108.483	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.787	88	108502	36.660	ng	98
3) Pyridine	3.522	79	293912	39.958	ng	98
4) n-Nitrosodimethylamine	3.487	42	190306	44.441	ng	96
6) Aniline	6.551	93	328800	40.838	ng	92
8) 2-Chlorophenol	6.675	128	323058	54.332	ng	99
9) Benzaldehyde	6.440	77	25756	4.791	ng	99
10) Phenol	6.534	94	439081	50.653	ng	99
11) bis(2-Chloroethyl)ether	6.628	93	324959	47.579	ng	99
12) 1,3-Dichlorobenzene	6.834	146	348562	44.454	ng	99
13) 1,4-Dichlorobenzene	6.910	146	355192	45.151	ng	100
14) 1,2-Dichlorobenzene	7.063	146	339112	45.816	ng	98
15) Benzyl Alcohol	7.034	79	341041	52.678	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.169	45	479204	48.400	ng	99
17) 2-Methylphenol	7.146	107	274650	50.669	ng	99
18) Hexachloroethane	7.404	117	127908	51.392	ng	99
19) n-Nitroso-di-n-propyla...	7.310	70	264857	51.557	ng	99
20) 3+4-Methylphenols	7.298	107	345199	49.926	ng	97
22) Acetophenone	7.304	105	465602	45.826	ng	99
24) Nitrobenzene	7.475	77	398115	47.881	ng	96
25) Isophorone	7.716	82	711016	50.373	ng	100
26) 2-Nitrophenol	7.793	139	160916	59.767	ng	98
27) 2,4-Dimethylphenol	7.828	122	302150	72.223	ng	98
28) bis(2-Chloroethoxy)met...	7.922	93	399132	48.841	ng	100
29) 2,4-Dichlorophenol	8.028	162	279528	54.132	ng	99
30) 1,2,4-Trichlorobenzene	8.116	180	304642	45.748	ng	99
31) Naphthalene	8.198	128	925097	45.753	ng	100
32) Benzoic acid	7.945	122	198469	53.987	ng	98
33) 4-Chloroaniline	8.240	127	181345	26.026	ng	99
34) Hexachlorobutadiene	8.310	225	194329	44.908	ng	99
35) Caprolactam	8.610	113	87995m	55.549	ng	
36) 4-Chloro-3-methylphenol	8.722	107	308098	52.023	ng	99
37) 2-Methylnaphthalene	8.887	142	591930	46.699	ng	98
38) 1-Methylnaphthalene	8.987	142	552231	44.575	ng	100
40) 1,2,4,5-Tetrachloroben...	9.051	216	299159	46.231	ng	98
41) Hexachlorocyclopentadiene	9.040	237	410655	197.633	ng	98
43) 2,4,6-Trichlorophenol	9.163	196	209677	57.950	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.204	196	216170	56.997	ng	98
46) 1,1'-Biphenyl	9.351	154	745616	47.217	ng	100
47) 2-Chloronaphthalene	9.375	162	586615	46.564	ng	98
48) 2-Nitroaniline	9.469	65	207561	55.009	ng	97
49) Acenaphthylene	9.792	152	894874	50.411	ng	100
50) Dimethylphthalate	9.651	163	694072	55.074	ng	100
51) 2,6-Dinitrotoluene	9.710	165	153547	55.959	ng	92
52) Acenaphthene	9.963	154	650721	55.802	ng	100
53) 3-Nitroaniline	9.881	138	106972	39.731	ng	95
54) 2,4-Dinitrophenol	9.986	184	169255	109.329	ng	# 48
55) Dibenzofuran	10.134	168	829838	48.760	ng	99
56) 4-Nitrophenol	10.039	139	257665	109.890	ng	98
57) 2,4-Dinitrotoluene	10.116	165	200941	57.411	ng	95
58) Fluorene	10.475	166	649223	48.353	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.251	232	187011	61.523	ng	99
60) Diethylphthalate	10.351	149	658383	55.478	ng	100
61) 4-Chlorophenyl-phenyle...	10.469	204	320467	48.634	ng	98
62) 4-Nitroaniline	10.498	138	144318	51.715	ng	97
63) Azobenzene	10.628	77	723378	50.141	ng	100
65) 4,6-Dinitro-2-methylph...	10.522	198	107485	63.201	ng	93
66) n-Nitrosodiphenylamine	10.586	169	563456	48.161	ng	99
67) 4-Bromophenyl-phenylether	10.957	248	203565	48.498	ng	99
68) Hexachlorobenzene	11.022	284	218339	48.066	ng	99
69) Atrazine	11.116	200	180535	70.747	ng	99
70) Pentachlorophenol	11.216	266	274506	102.930	ng	99
71) Phenanthrene	11.439	178	969096	48.546	ng	100
72) Anthracene	11.492	178	967869	49.738	ng	100
73) Carbazole	11.645	167	822196	47.232	ng	100
74) Di-n-butylphthalate	11.975	149	914192	60.594	ng	100
75) Fluoranthene	12.628	202	937525	48.394	ng	99
77) Benzidine	12.745	184	105216	32.172	ng	99
78) Pyrene	12.851	202	914365	50.226	ng	100
80) Butylbenzylphthalate	13.475	149	239346	62.503	ng	100
81) Benzo(a)anthracene	14.039	228	628151	48.916	ng	100
82) 3,3'-Dichlorobenzidine	14.004	252	149055	49.854	ng	100
83) Chrysene	14.074	228	569951	49.203	ng	99
84) Bis(2-ethylhexyl)phtha...	14.033	149	275070	63.368	ng	98
85) Di-n-octyl phthalate	14.645	149	522438	58.475	ng	98
87) Indeno(1,2,3-cd)pyrene	17.004	276	689061	53.881	ng	100
88) Benzo(b)fluoranthene	15.092	252	575818	52.630	ng	99
89) Benzo(k)fluoranthene	15.121	252	556895	56.300	ng	99
90) Benzo(a)pyrene	15.457	252	505638	54.448	ng	100
91) Dibenzo(a,h)anthracene	17.027	278	577395	55.085	ng	98
92) Benzo(g,h,i)perylene	17.457	276	567774	52.368	ng	98

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

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