

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF052324\
 Data File : BF138020.D
 Acq On : 23 May 2024 11:21
 Operator : CG\JU
 Sample : PB161055BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB161055BS

Quant Time: May 23 11:56:00 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF051424.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 15 04:12:16 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.034	152	159557	20.000	ng	-0.01
21) Naphthalene-d8	8.322	136	634354	20.000	ng	0.00
39) Acenaphthene-d10	10.086	164	360690	20.000	ng	0.00
64) Phenanthrene-d10	11.580	188	631196	20.000	ng	0.00
76) Chrysene-d12	14.233	240	354682	20.000	ng	0.00
86) Perylene-d12	15.798	264	306386	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.675	112	1085622	115.575	ng	0.01
7) Phenol-d6	6.657	99	1347805	116.482	ng	0.00
23) Nitrobenzene-d5	7.604	82	827658	76.784	ng	0.00
42) 2,4,6-Tribromophenol	10.880	330	472751	129.990	ng	0.00
45) 2-Fluorobiphenyl	9.398	172	1776940	75.887	ng	0.00
79) Terphenyl-d14	13.168	244	1955763	81.267	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.040	88	126984	31.206	ng	97
3) Pyridine	3.787	79	317553	34.780	ng	100
4) n-Nitrosodimethylamine	3.751	42	164568	40.857	ng	94
6) Aniline	6.698	93	323013m	31.211	ng	
8) 2-Chlorophenol	6.822	128	439953	43.280	ng	99
9) Benzaldehyde	6.592	77	363366	58.044	ng	99
10) Phenol	6.675	94	508361	43.527	ng	93
11) bis(2-Chloroethyl)ether	6.769	93	381838	42.580	ng	98
12) 1,3-Dichlorobenzene	6.975	146	460924	39.190	ng	98
13) 1,4-Dichlorobenzene	7.051	146	467514	39.601	ng	98
14) 1,2-Dichlorobenzene	7.204	146	450340	40.531	ng	99
15) Benzyl Alcohol	7.175	79	356788	46.177	ng	98
16) 2,2'-oxybis(1-Chloropr...	7.298	45	475066	43.097	ng	98
17) 2-Methylphenol	7.281	107	352365	43.466	ng	97
18) Hexachloroethane	7.551	117	159891	40.024	ng	96
19) n-Nitroso-di-n-propyla...	7.445	70	282860	43.512	ng	98
20) 3+4-Methylphenols	7.428	107	444998	41.901	ng	93
22) Acetophenone	7.445	105	584360	41.121	ng	# 98
24) Nitrobenzene	7.622	77	426883	39.038	ng	99
25) Isophorone	7.857	82	785882	42.008	ng	98
26) 2-Nitrophenol	7.933	139	242445	44.125	ng	99
27) 2,4-Dimethylphenol	7.957	122	315593	44.613	ng	98
28) bis(2-Chloroethoxy)met...	8.057	93	476803	42.052	ng	99
29) 2,4-Dichlorophenol	8.175	162	381970	42.939	ng	97
30) 1,2,4-Trichlorobenzene	8.257	180	394009	39.670	ng	99
31) Naphthalene	8.345	128	1264611	39.649	ng	100
32) Benzoic acid	8.086	122	285927	47.815	ng	99
33) 4-Chloroaniline	8.386	127	217606	20.812	ng	98
34) Hexachlorobutadiene	8.451	225	239465	38.091	ng	99
35) Caprolactam	8.769	113	120129m	46.612	ng	
36) 4-Chloro-3-methylphenol	8.863	107	393917	43.531	ng	97
37) 2-Methylnaphthalene	9.039	142	846609	41.293	ng	99
38) 1-Methylnaphthalene	9.139	142	788154	39.103	ng	99
40) 1,2,4,5-Tetrachloroben...	9.204	216	405322	41.335	ng	99
41) Hexachlorocyclopentadiene	9.186	237	405833	108.921	ng	99
43) 2,4,6-Trichlorophenol	9.310	196	279477	43.469	ng	99

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.357	196	298965	42.498	ng	99
46) 1,1'-Biphenyl	9.504	154	1037740	41.020	ng	99
47) 2-Chloronaphthalene	9.533	162	806371	39.930	ng	100
48) 2-Nitroaniline	9.627	65	227759	44.042	ng	90
49) Acenaphthylene	9.951	152	1293817	44.557	ng	99
50) Dimethylphthalate	9.804	163	980547	43.352	ng	100
51) 2,6-Dinitrotoluene	9.869	165	222315	42.826	ng	88
52) Acenaphthene	10.122	154	869950	45.329	ng	99
53) 3-Nitroaniline	10.039	138	164799	32.473	ng	97
54) 2,4-Dinitrophenol	10.145	184	238454	84.588	ng	# 85
55) Dibenzofuran	10.292	168	1162551	41.622	ng	99
56) 4-Nitrophenol	10.192	139	341312	93.639	ng	98
57) 2,4-Dinitrotoluene	10.275	165	299004	44.554	ng	97
58) Fluorene	10.639	166	924382	41.842	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.410	232	251192	45.453	ng	99
60) Diethylphthalate	10.504	149	913024	43.222	ng	99
61) 4-Chlorophenyl-phenyle...	10.627	204	455912	41.846	ng	94
62) 4-Nitroaniline	10.657	138	202663	42.104	ng	100
63) Azobenzene	10.786	77	827285	43.056	ng	99
65) 4,6-Dinitro-2-methylph...	10.680	198	170587	46.806	ng	90
66) n-Nitrosodiphenylamine	10.745	169	795314	41.087	ng	99
67) 4-Bromophenyl-phenylether	11.121	248	289784	41.878	ng	91
68) Hexachlorobenzene	11.186	284	312178	40.813	ng	95
69) Atrazine	11.269	200	273898	51.819	ng	99
70) Pentachlorophenol	11.380	266	373973	83.467	ng	99
71) Phenanthrene	11.610	178	1292715	41.560	ng	99
72) Anthracene	11.663	178	1303690	42.243	ng	99
73) Carbazole	11.816	167	1171864	41.567	ng	99
74) Di-n-butylphthalate	12.127	149	1378707	44.067	ng	99
75) Fluoranthene	12.804	202	1305974	41.365	ng	99
77) Benzidine	12.916	184	93390	20.148	ng	98
78) Pyrene	13.033	202	1300504	41.922	ng	99
80) Butylbenzylphthalate	13.633	149	474625	50.724	ng	95
81) Benzo(a)anthracene	14.221	228	985696	43.927	ng	99
82) 3,3'-Dichlorobenzidine	14.180	252	212780	34.217	ng	99
83) Chrysene	14.262	228	919960	42.868	ng	99
84) Bis(2-ethylhexyl)phtha...	14.186	149	578936	49.219	ng	97
85) Di-n-octyl phthalate	14.815	149	800457	50.374	ng	99
87) Indeno(1,2,3-cd)pyrene	17.433	276	904918	45.971	ng	99
88) Benzo(b)fluoranthene	15.327	252	786048	44.310	ng	99
89) Benzo(k)fluoranthene	15.362	252	809143	46.428	ng	99
90) Benzo(a)pyrene	15.733	252	720487	47.563	ng	99
91) Dibenzo(a,h)anthracene	17.456	278	735752	45.239	ng	99
92) Benzo(g,h,i)perylene	17.933	276	701985	41.088	ng	99

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

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