

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061523\
 Data File : BF133799.D
 Acq On : 16 Jun 2023 04:13
 Operator : CG\JU
 Sample : 03221-02MSD
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RT-4194MSD

Quant Time: Jun 16 05:33:40 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 15 15:12:57 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.834	152	108696	20.000	ng	0.00
21) Naphthalene-d8	8.116	136	392243	20.000	ng	0.00
39) Acenaphthene-d10	9.869	164	166502	20.000	ng	0.00
64) Phenanthrene-d10	11.357	188	251981	20.000	ng	0.00
76) Chrysene-d12	14.004	240	206800	20.000	ng	0.00
86) Perylene-d12	15.463	264	235375	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.446	112	420210	67.302	ng	0.00
7) Phenol-d6	6.463	99	503903	62.002	ng	0.00
23) Nitrobenzene-d5	7.398	82	341154	47.373	ng	0.00
42) 2,4,6-Tribromophenol	10.657	330	108961	51.699	ng	0.00
45) 2-Fluorobiphenyl	9.186	172	620676	52.990	ng	0.00
79) Terphenyl-d14	12.939	244	493034	36.890	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.610	88	108148	39.399	ng	97
3) Pyridine	3.357	79	285674	35.706	ng	96
4) n-Nitrosodimethylamine	3.316	42	196668	44.470	ng	97
6) Aniline	6.493	93	400720	39.147	ng	100
8) 2-Chlorophenol	6.616	128	317963	48.324	ng	96
9) Benzaldehyde	6.381	77	181374	33.638	ng	96
10) Phenol	6.475	94	395147	46.563	ng	99
11) bis(2-Chloroethyl)ether	6.569	93	291793	46.315	ng	96
12) 1,3-Dichlorobenzene	6.775	146	342362	48.699	ng	99
13) 1,4-Dichlorobenzene	6.851	146	346775	48.803	ng	97
14) 1,2-Dichlorobenzene	6.998	146	330018	51.400	ng	98
15) Benzyl Alcohol	6.975	79	289067	45.540	ng	97
16) 2,2'-oxybis(1-Chloropr...	7.110	45	523729	49.039	ng	100
17) 2-Methylphenol	7.087	107	255584	43.824	ng	100
18) Hexachloroethane	7.340	117	128739	52.923	ng	94
19) n-Nitroso-di-n-propyla...	7.245	70	223110	43.147	ng	97
20) 3+4-Methylphenols	7.240	107	304842	40.181	ng	# 89
22) Acetophenone	7.245	105	432247	48.047	ng	# 97
24) Nitrobenzene	7.416	77	345198	47.604	ng	99
25) Isophorone	7.657	82	610104	46.146	ng	98
26) 2-Nitrophenol	7.734	139	173419	53.288	ng	97
27) 2,4-Dimethylphenol	7.769	122	271829	48.461	ng	99
28) bis(2-Chloroethoxy)met...	7.863	93	352297	46.104	ng	97
29) 2,4-Dichlorophenol	7.975	162	250930	45.644	ng	98
30) 1,2,4-Trichlorobenzene	8.057	180	274665	47.887	ng	98
31) Naphthalene	8.139	128	856466	44.818	ng	99
32) Benzoic acid	7.875	122	154690	43.727	ng	95
33) 4-Chloroaniline	8.187	127	267993	32.798	ng	98
34) Hexachlorobutadiene	8.251	225	173781	46.276	ng	97
35) Caprolactam	8.563	113	71426m	38.611	ng	
36) 4-Chloro-3-methylphenol	8.663	107	261714	41.301	ng	96
37) 2-Methylnaphthalene	8.828	142	562377	42.354	ng	100
38) 1-Methylnaphthalene	8.928	142	519593	42.526	ng	100
40) 1,2,4,5-Tetrachloroben...	8.992	216	262182	51.796	ng	99
41) Hexachlorocyclopentadiene	8.981	237	239446	106.151	ng	97
43) 2,4,6-Trichlorophenol	9.104	196	167212	49.434	ng	98

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44) 2,4,5-Trichlorophenol	9.145	196	168048	42.704	ng	98
46) 1,1'-Biphenyl	9.292	154	683860	50.926	ng	99
47) 2-Chloronaphthalene	9.316	162	493820	51.266	ng	99
48) 2-Nitroaniline	9.416	65	165328	46.930	ng	98
49) Acenaphthylene	9.733	152	780946	46.580	ng	100
50) Dimethylphthalate	9.592	163	571375	45.975	ng	99
51) 2,6-Dinitrotoluene	9.657	165	119589	46.619	ng	# 84
52) Acenaphthene	9.904	154	459986	46.873	ng	99
53) 3-Nitroaniline	9.828	138	116578	37.344	ng	94
54) 2,4-Dinitrophenol	9.933	184	94000	75.346	ng	99
55) Dibenzofuran	10.075	168	676228	45.159	ng	97
56) 4-Nitrophenol	9.986	139	164619	78.135	ng	84
57) 2,4-Dinitrotoluene	10.057	165	143692	44.044	ng	97
58) Fluorene	10.422	166	497884	43.478	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.192	232	125437	42.576	ng	99
60) Diethylphthalate	10.292	149	536392	42.407	ng	100
61) 4-Chlorophenyl-phenyle...	10.410	204	243448	43.064	ng	99
62) 4-Nitroaniline	10.439	138	106453	34.779	ng	96
63) Azobenzene	10.569	77	513656	42.897	ng	99
65) 4,6-Dinitro-2-methylph...	10.469	198	62365	51.794	ng	100
66) n-Nitrosodiphenylamine	10.528	169	425489	50.270	ng	99
67) 4-Bromophenyl-phenylether	10.904	248	134356	47.772	ng	92
68) Hexachlorobenzene	10.963	284	143481	48.514	ng	97
69) Atrazine	11.057	200	120788	49.184	ng	99
70) Pentachlorophenol	11.163	266	147575	83.330	ng	99
71) Phenanthrene	11.380	178	634646	49.643	ng	99
72) Anthracene	11.433	178	614025	46.063	ng	99
73) Carbazole	11.586	167	565403	45.541	ng	99
74) Di-n-butylphthalate	11.916	149	641000	40.274	ng	99
75) Fluoranthene	12.574	202	764447	56.309	ng	100
77) Benzidine	12.698	184	355621	51.247	ng	99
78) Pyrene	12.804	202	807699	41.927	ng	100
80) Butylbenzylphthalate	13.421	149	289865	35.642	ng	99
81) Benzo(a)anthracene	13.992	228	764277	53.429	ng	98
82) 3,3'-Dichlorobenzidine	13.957	252	241464	50.774	ng	98
83) Chrysene	14.033	228	729826	53.944	ng	99
84) Bis(2-ethylhexyl)phtha...	13.980	149	410871	41.031	ng	99
85) Di-n-octyl phthalate	14.598	149	786673	55.283	ng	97
87) Indeno(1,2,3-cd)pyrene	16.927	276	689929	42.606	ng	99
88) Benzo(b)fluoranthene	15.045	252	921944	64.480	ng	99
89) Benzo(k)fluoranthene	15.074	252	735114	52.780	ng	98
90) Benzo(a)pyrene	15.404	252	713936	53.802	ng	99
91) Dibenzo(a,h)anthracene	16.945	278	534609	39.878	ng	99
92) Benzo(g,h,i)perylene	17.374	276	541566	40.744	ng	98

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

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