

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF112222\
 Data File : BF131348.D
 Acq On : 23 Nov 2022 00:57
 Operator : CG\JU
 Sample : N5740-01MSD
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MH-12MSD

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 11/23/2022
 Supervised By : Jagrut Upadhyay 11/23/2022

Quant Time: Nov 23 04:22:06 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF112122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 22 01:17:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.963	152	89896	20.000	ng	0.00
21) Naphthalene-d8	8.257	136	324650	20.000	ng	# 0.00
39) Acenaphthene-d10	10.022	164	169117	20.000	ng	0.00
64) Phenanthrene-d10	11.516	188	247646	20.000	ng	0.00
76) Chrysene-d12	14.174	240	182361	20.000	ng	0.00
86) Perylene-d12	15.715	264	198108	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.581	112	591002	125.123	ng	0.01
7) Phenol-d6	6.581	99	738540	122.549	ng	0.00
23) Nitrobenzene-d5	7.534	82	473893	73.582	ng	0.00
42) 2,4,6-Tribromophenol	10.816	330	175739	123.517	ng	0.00
45) 2-Fluorobiphenyl	9.334	172	819553	70.502	ng	0.00
79) Terphenyl-d14	13.110	244	628551	56.363	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.834	88	92965	41.284	ng	98
3) Pyridine	3.593	79	248884	39.804	ng	95
4) n-Nitrosodimethylamine	3.557	42	166878	45.257	ng	98
6) Aniline	6.628	93	208932m	26.454	ng	
8) 2-Chlorophenol	6.745	128	271294	50.698	ng	99
9) Benzaldehyde	6.516	77	153103	41.772	ng	97
10) Phenol	6.598	94	324700	48.595	ng	96
11) bis(2-Chloroethyl)ether	6.698	93	262163	46.830	ng	99
12) 1,3-Dichlorobenzene	6.904	146	294212	46.132	ng	97
13) 1,4-Dichlorobenzene	6.981	146	297574	45.919	ng	97
14) 1,2-Dichlorobenzene	7.134	146	282331	47.316	ng	99
15) Benzyl Alcohol	7.104	79	241665	48.906	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.240	45	476188	44.478	ng	99
17) 2-Methylphenol	7.210	107	218361	47.877	ng	98
18) Hexachloroethane	7.481	117	113392	48.914	ng	96
19) n-Nitroso-di-n-propyla...	7.381	70	195670	44.352	ng	97
20) 3+4-Methylphenols	7.363	107	272303	46.760	ng	97
22) Acetophenone	7.381	105	378940	45.856	ng	# 95
24) Nitrobenzene	7.551	77	300803	44.828	ng	99
25) Isophorone	7.792	82	542469	47.211	ng	97
26) 2-Nitrophenol	7.869	139	131544	58.473	ng	92
27) 2,4-Dimethylphenol	7.898	122	238325	53.503	ng	99
28) bis(2-Chloroethoxy)met...	7.998	93	318514	48.806	ng	98
29) 2,4-Dichlorophenol	8.104	162	226911	46.963	ng	97
30) 1,2,4-Trichlorobenzene	8.192	180	260994	43.604	ng	99
31) Naphthalene	8.281	128	740947	42.828	ng	99
32) Benzoic acid	7.992	122	115596m	43.308	ng	
33) 4-Chloroaniline	8.322	127	51647	7.567	ng	95
34) Hexachlorobutadiene	8.392	225	168206	43.399	ng	98
35) Caprolactam	8.698	113	64854m	51.080	ng	
36) 4-Chloro-3-methylphenol	8.798	107	236276	47.069	ng	93
37) 2-Methylnaphthalene	8.969	142	500529	42.700	ng	100
38) 1-Methylnaphthalene	9.075	142	466121	41.007	ng	97
40) 1,2,4,5-Tetrachloroben...	9.134	216	262108	47.760	ng	100
41) Hexachlorocyclopentadiene	9.122	237	265801	107.737	ng	98
43) 2,4,6-Trichlorophenol	9.245	196	165352	53.111	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.281	196	170556	48.718	ng	95
46) 1,1'-Biphenyl	9.439	154	608499	47.150	ng	99
47) 2-Chloronaphthalene	9.463	162	477019	46.076	ng	99
48) 2-Nitroaniline	9.557	65	158408	50.394	ng	97
49) Acenaphthylene	9.881	152	707989	44.861	ng	100
50) Dimethylphthalate	9.739	163	553900	47.008	ng	99
51) 2,6-Dinitrotoluene	9.798	165	117357	49.803	ng	96
52) Acenaphthene	10.057	154	462427	46.793	ng	98
53) 3-Nitroaniline	9.969	138	54162	22.368	ng	96
54) 2,4-Dinitrophenol	10.075	184	70477	69.673	ng	# 80
55) Dibenzofuran	10.228	168	616266	43.346	ng	99
56) 4-Nitrophenol	10.122	139	155881	92.760	ng	96
57) 2,4-Dinitrotoluene	10.204	165	148176	49.977	ng	97
58) Fluorene	10.575	166	468402	42.214	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.339	232	131757	45.192	ng	100
60) Diethylphthalate	10.439	149	502048	45.124	ng	99
61) 4-Chlorophenyl-phenyle...	10.563	204	250536	44.045	ng	99
62) 4-Nitroaniline	10.586	138	93442	38.776	ng	97
63) Azobenzene	10.722	77	501874	41.952	ng	99
65) 4,6-Dinitro-2-methylph...	10.610	198	49390	46.828	ng	87
66) n-Nitrosodiphenylamine	10.680	169	408530	51.771	ng	99
67) 4-Bromophenyl-phenylether	11.057	248	152762	50.330	ng	100
68) Hexachlorobenzene	11.116	284	159982	49.662	ng	99
69) Atrazine	11.210	200	139973	57.611	ng	99
70) Pentachlorophenol	11.310	266	152075	81.168	ng	99
71) Phenanthrene	11.539	178	605319	45.228	ng	100
72) Anthracene	11.592	178	611158	46.465	ng	99
73) Carbazole	11.745	167	504584	44.973	ng	99
74) Di-n-butylphthalate	12.075	149	650287	49.187	ng	100
75) Fluoranthene	12.733	202	571672	40.536	ng	100
77) Benzidine	12.857	184	219671	65.215	ng	97
78) Pyrene	12.969	202	573385	35.637	ng	99
80) Butylbenzylphthalate	13.586	149	222571	41.712	ng	97
81) Benzo(a)anthracene	14.163	228	549436	43.955	ng	100
82) 3,3'-Dichlorobenzidine	14.121	252	105412	31.780	ng	98
83) Chrysene	14.198	228	533425	43.496	ng	99
84) Bis(2-ethylhexyl)phtha...	14.145	149	309479	43.466	ng	100
85) Di-n-octyl phthalate	14.774	149	570832	53.904	ng	97
87) Indeno(1,2,3-cd)pyrene	17.292	276	541931	40.729	ng	99
88) Benzo(b)fluoranthene	15.257	252	593149	45.070	ng	99
89) Benzo(k)fluoranthene	15.292	252	601829	44.499	ng	97
90) Benzo(a)pyrene	15.651	252	535209	49.562	ng	99
91) Dibenzo(a,h)anthracene	17.315	278	464314	42.267	ng	98
92) Benzo(g,h,i)perylene	17.768	276	422840	38.566	ng	98

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

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