

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072921\
 Data File : BF124855.D
 Acq On : 29 Jul 2021 17:14
 Operator : JU/CG
 Sample : M3176-02MS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 18-B7-TP-(9.5-12)MS

Manual Integrations
 APPROVED

mohammad
 7/30/2021 12:00:55 PM

Quant Time: Jul 29 18:24:25 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF072221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 22 15:28:57 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.875	152	7289	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	29346	20.000	ng	# 0.00
39) Acenaphthene-d10	9.916	164	16926	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	32063	20.000	ng	0.00
76) Chrysene-d12	14.051	240	21852	20.000	ng	# 0.01
86) Perylene-d12	15.521	264	16939	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	54977	127.706	ng	0.03
7) Phenol-d6	6.510	99	62055	114.982	ng	0.01
23) Nitrobenzene-d5	7.445	82	48601	98.569	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	23143	159.249	ng	0.00
45) 2-Fluorobiphenyl	9.233	172	103322	89.657	ng	0.00
79) Terphenyl-d14	12.992	244	130540	102.400	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.863	88	6363	42.284	ng	# 100
3) Pyridine	3.551	79	12398	34.804	ng	# 90
4) n-Nitrosodimethylamine	3.498	42	12523	44.453	ng	# 80
6) Aniline	6.533	93	11861	21.385	ng	# 88
8) 2-Chlorophenol	6.663	128	23340	48.069	ng	92
9) Benzaldehyde	6.428	77	9149	31.486	ng	93
10) Phenol	6.522	94	23102	44.700	ng	92
11) bis(2-Chloroethyl)ether	6.610	93	23446	57.216	ng	98
12) 1,3-Dichlorobenzene	6.816	146	22768	43.143	ng	97
13) 1,4-Dichlorobenzene	6.892	146	23975	45.411	ng	98
14) 1,2-Dichlorobenzene	7.045	146	22710	44.520	ng	95
15) Benzyl Alcohol	7.022	79	19142	53.936	ng	91
16) 2,2'-oxybis(1-Chloropr...	7.151	45	35106	50.900	ng	94
17) 2-Methylphenol	7.128	107	17697	50.029	ng	95
18) Hexachloroethane	7.386	117	9910	49.523	ng	94
19) n-Nitroso-di-n-propyla...	7.292	70	16518	57.317	ng	92
20) 3+4-Methylphenols	7.281	107	23120	49.469	ng	# 77
22) Acetophenone	7.286	105	33744	49.620	ng	# 92
24) Nitrobenzene	7.463	77	24716	51.129	ng	98
25) Isophorone	7.698	82	44404	50.931	ng	# 92
26) 2-Nitrophenol	7.775	139	13022	48.637	ng	# 90
27) 2,4-Dimethylphenol	7.810	122	20622	50.056	ng	93
28) bis(2-Chloroethoxy)met...	7.904	93	28501	56.254	ng	# 96
29) 2,4-Dichlorophenol	8.016	162	20334	46.603	ng	97
30) 1,2,4-Trichlorobenzene	8.098	180	21463	44.882	ng	99
31) Naphthalene	8.180	128	70310	45.910	ng	97
32) Benzoic acid	7.945	122	11327	35.521	ng	87
33) 4-Chloroaniline	8.222	127	5991	10.403	ng	# 87
34) Hexachlorobutadiene	8.292	225	13544	47.380	ng	97
35) Caprolactam	8.622	113	5633m	49.672	ng	
36) 4-Chloro-3-methylphenol	8.716	107	22983	55.224	ng	93
37) 2-Methylnaphthalene	8.869	142	48906	47.794	ng	98
38) 1-Methylnaphthalene	8.969	142	44799	46.213	ng	95
40) 1,2,4,5-Tetrachloroben...	9.039	216	22077	46.748	ng	98
41) Hexachlorocyclopentadiene	9.022	237	23273	83.295	ng	94
43) 2,4,6-Trichlorophenol	9.151	196	16020	47.785	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.192	196	16856	48.967	ng	96
46) 1,1'-Biphenyl	9.339	154	58831	46.564	ng	99
47) 2-Chloronaphthalene	9.363	162	47189	47.181	ng	97
48) 2-Nitroaniline	9.463	65	14346	54.778	ng	96
49) Acenaphthylene	9.780	152	73392	47.583	ng	99
50) Dimethylphthalate	9.645	163	62089	53.274	ng	# 97
51) 2,6-Dinitrotoluene	9.704	165	12841	49.872	ng	89
52) Acenaphthene	9.951	154	45117	44.125	ng	97
53) 3-Nitroaniline	9.874	138	7995	29.515	ng	93
54) 2,4-Dinitrophenol	9.980	184	10757	72.249	ng	91
55) Dibenzofuran	10.127	168	69622	47.651	ng	98
56) 4-Nitrophenol	10.039	139	19034	88.981	ng	88
57) 2,4-Dinitrotoluene	10.110	165	17554	50.239	ng	# 95
58) Fluorene	10.469	166	54736	48.835	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.239	232	13926	51.225	ng	# 96
60) Diethylphthalate	10.339	149	66333	54.733	ng	98
61) 4-Chlorophenyl-phenyle...	10.457	204	27624	50.983	ng	96
62) 4-Nitroaniline	10.492	138	12808	47.785	ng	94
63) Azobenzene	10.616	77	57367	53.504	ng	93
65) 4,6-Dinitro-2-methylph...	10.521	198	8071	38.869	ng	89
66) n-Nitrosodiphenylamine	10.580	169	48257	46.444	ng	96
67) 4-Bromophenyl-phenylether	10.945	248	17442	51.293	ng	# 88
68) Hexachlorobenzene	11.016	284	18873	53.441	ng	# 88
69) Atrazine	11.110	200	17075	54.051	ng	94
70) Pentachlorophenol	11.210	266	19379	88.245	ng	91
71) Phenanthrene	11.433	178	82529	47.857	ng	99
72) Anthracene	11.486	178	84307	48.902	ng	99
73) Carbazole	11.639	167	74061	45.829	ng	98
74) Di-n-butylphthalate	11.963	149	114110	53.013	ng	99
75) Fluoranthene	12.621	202	84451	47.989	ng	96
77) Benzidine	12.733	184	14848	24.096	ng	98
78) Pyrene	12.851	202	85122	49.205	ng	99
80) Butylbenzylphthalate	13.462	149	44707	53.956	ng	98
81) Benzo(a)anthracene	14.033	228	66166	46.322	ng	98
82) 3,3'-Dichlorobenzidine	13.998	252	13409	33.826	ng	98
83) Chrysene	14.074	228	62145	45.159	ng	98
84) Bis(2-ethylhexyl)phtha...	14.015	149	69642	57.061	ng	99
85) Di-n-octyl phthalate	14.633	149	105238	53.137	ng	100
87) Indeno(1,2,3-cd)pyrene	17.015	276	52046	53.220	ng	94
88) Benzo(b)fluoranthene	15.092	252	53654m	44.988	ng	
89) Benzo(k)fluoranthene	15.121	252	50834	46.649	ng	99
90) Benzo(a)pyrene	15.462	252	48527	48.709	ng	97
91) Dibenzo(a,h)anthracene	17.033	278	45165	52.756	ng	94
92) Benzo(g,h,i)perylene	17.468	276	42323	52.133	ng	# 88

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

