

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072121\
 Data File : BF124669.D
 Acq On : 21 Jul 2021 19:11
 Operator : JU/CG
 Sample : M3076-04MSD
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20210587-AMENDED-COM-1-SPLP

Manual Integrations
 APPROVED

mohammad
 7/22/2021 1:29:02 PM

Quant Time: Jul 22 01:29:29 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF070721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 21 16:08:04 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.886	152	5934	20.000	ng	0.00
21) Naphthalene-d8	8.175	136	23911	20.000	ng	0.00
39) Acenaphthene-d10	9.933	164	13600	20.000	ng	0.00
64) Phenanthrene-d10	11.421	188	24361	20.000	ng	0.00
76) Chrysene-d12	14.062	240	15286	20.000	ng	# 0.00
86) Perylene-d12	15.539	264	14871	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	33787	98.649	ng	0.00
7) Phenol-d6	6.510	99	28064	67.357	ng	-0.01
23) Nitrobenzene-d5	7.457	82	40540	107.029	ng	0.00
42) 2,4,6-Tribromophenol	10.727	330	18157	171.562	ng	0.00
45) 2-Fluorobiphenyl	9.251	172	97718	104.777	ng	0.00
79) Terphenyl-d14	13.010	244	98633	108.351	ng	0.00
Target Compounds						
2) 1,4-Dioxane	2.704	88	3678	31.979	ng	# 100
3) Pyridine	3.463	79	6284	21.871	ng	# 88
4) n-Nitrosodimethylamine	3.404	42	7328m	31.379	ng	
6) Aniline	6.551	93	19425	44.248	ng	97
8) 2-Chlorophenol	6.675	128	18838	49.061	ng	97
9) Benzaldehyde	6.439	77	13464	55.082	ng	91
10) Phenol	6.522	94	10568	25.980	ng	95
11) bis(2-Chloroethyl)ether	6.622	93	18979	60.181	ng	98
12) 1,3-Dichlorobenzene	6.833	146	20870	49.660	ng	96
13) 1,4-Dichlorobenzene	6.910	146	20418	47.567	ng	96
14) 1,2-Dichlorobenzene	7.063	146	20508	50.273	ng	98
15) Benzyl Alcohol	7.028	79	13022	45.169	ng	96
16) 2,2'-oxybis(1-Chloropr...	7.163	45	32162	59.169	ng	98
17) 2-Methylphenol	7.139	107	13088	46.567	ng	97
18) Hexachloroethane	7.404	117	8012	49.566	ng	93
19) n-Nitroso-di-n-propyla...	7.304	70	13908	59.373	ng	96
20) 3+4-Methylphenols	7.292	107	15755	42.577	ng	88
22) Acetophenone	7.298	105	30568	56.331	ng	# 95
24) Nitrobenzene	7.475	77	19041	50.697	ng	96
25) Isophorone	7.716	82	37603	54.079	ng	96
26) 2-Nitrophenol	7.786	139	11862	58.548	ng	92
27) 2,4-Dimethylphenol	7.822	122	19717	58.730	ng	98
28) bis(2-Chloroethoxy)met...	7.922	93	25484	62.830	ng	97
29) 2,4-Dichlorophenol	8.028	162	18258	52.725	ng	98
30) 1,2,4-Trichlorobenzene	8.116	180	19139	49.943	ng	98
31) Naphthalene	8.198	128	62887	50.850	ng	97
32) Benzoic acid	7.916	122	7370	31.004	ng	99
33) 4-Chloroaniline	8.239	127	19116	40.904	ng	98
34) Hexachlorobutadiene	8.310	225	10897	50.378	ng	98
35) Caprolactam	8.627	113	1282m	14.637	ng	
36) 4-Chloro-3-methylphenol	8.716	107	18217	53.200	ng	99
37) 2-Methylnaphthalene	8.886	142	44199	52.963	ng	99
38) 1-Methylnaphthalene	8.986	142	41816	53.354	ng	98
40) 1,2,4,5-Tetrachloroben...	9.057	216	19977	55.074	ng	94
41) Hexachlorocyclopentadiene	9.045	237	26538	117.121	ng	94
43) 2,4,6-Trichlorophenol	9.163	196	14839	56.666	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.198	196	14923	55.254	ng	97
46) 1,1'-Biphenyl	9.357	154	54293	52.563	ng	100
47) 2-Chloronaphthalene	9.380	162	42505	52.486	ng	99
48) 2-Nitroaniline	9.474	65	11548	59.298	ng	96
49) Acenaphthylene	9.798	152	68828	54.462	ng	99
50) Dimethylphthalate	9.657	163	54703	57.813	ng	97
51) 2,6-Dinitrotoluene	9.716	165	11324	60.708	ng	# 87
52) Acenaphthene	9.969	154	49473	61.223	ng	97
53) 3-Nitroaniline	9.886	138	8223	39.505	ng	93
54) 2,4-Dinitrophenol	9.992	184	12836	131.876	ng	93
55) Dibenzofuran	10.139	168	63007	53.031	ng	99
56) 4-Nitrophenol	10.039	139	10258	59.212	ng	97
57) 2,4-Dinitrotoluene	10.121	165	15788	61.824	ng	# 89
58) Fluorene	10.486	166	48445	52.714	ng	97
59) 2,3,4,6-Tetrachlorophenol	10.257	232	12173	58.360	ng	# 95
60) Diethylphthalate	10.357	149	58638	59.270	ng	99
61) 4-Chlorophenyl-phenyle...	10.474	204	23941	57.393	ng	97
62) 4-Nitroaniline	10.504	138	11399	54.337	ng	91
63) Azobenzene	10.633	77	47446	54.864	ng	99
65) 4,6-Dinitro-2-methylph...	10.527	198	8283	59.521	ng	96
66) n-Nitrosodiphenylamine	10.598	169	42693	53.821	ng	95
67) 4-Bromophenyl-phenylether	10.969	248	14641	58.122	ng	97
68) Hexachlorobenzene	11.027	284	15118	58.577	ng	98
69) Atrazine	11.121	200	12580	51.783	ng	96
70) Pentachlorophenol	11.221	266	18178	111.352	ng	96
71) Phenanthrene	11.451	178	71021	53.437	ng	99
72) Anthracene	11.498	178	72874	54.655	ng	99
73) Carbazole	11.651	167	61162	49.643	ng	100
74) Di-n-butylphthalate	11.980	149	96908	58.391	ng	100
75) Fluoranthene	12.633	202	67013	49.975	ng	99
77) Benzidine	12.757	184	28245	49.427	ng	97
78) Pyrene	12.863	202	65133	51.286	ng	97
80) Butylbenzylphthalate	13.480	149	34324	55.897	ng	99
81) Benzo(a)anthracene	14.051	228	52303	51.592	ng	97
82) 3,3'-Dichlorobenzidine	14.009	252	14302	53.859	ng	95
83) Chrysene	14.086	228	51103	52.621	ng	98
84) Bis(2-ethylhexyl)phtha...	14.039	149	53565	64.972	ng	100
85) Di-n-octyl phthalate	14.651	149	89116	74.702	ng	99
87) Indeno(1,2,3-cd)pyrene	17.045	276	58010	67.261	ng	96
88) Benzo(b)fluoranthene	15.109	252	55055m	51.579	ng	
89) Benzo(k)fluoranthene	15.139	252	47765	48.772	ng	98
90) Benzo(a)pyrene	15.480	252	50280	57.066	ng	98
91) Dibenzo(a,h)anthracene	17.062	278	49742	67.759	ng	98
92) Benzo(g,h,i)perylene	17.497	276	47544	66.335	ng	90

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

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