

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071921\
 Data File : BF124613.D
 Acq On : 19 Jul 2021 14:44
 Operator : JU/CG
 Sample : PB137725BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB137725BS

Manual Integrations
 APPROVED

mohammad
 7/20/2021 11:37:48 AM

Quant Time: Jul 19 15:12:40 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF070721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 19 11:33:09 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.916	152	11089	20.00	ng	0.00
21) Naphthalene-d8	8.204	136	44643	20.00	ng	0.00
39) Acenaphthene-d10	9.963	164	25025	20.00	ng	0.00
64) Phenanthrene-d10	11.451	188	43329	20.00	ng	0.00
76) Chrysene-d12	14.092	240	28035	20.00	ng	0.00
86) Perylene-d12	15.580	264	19717	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.563	112	84554	132.11	ng	0.03
7) Phenol-d6	6.551	99	99096	127.27	ng	0.00
23) Nitrobenzene-d5	7.487	82	64315	90.94	ng	0.00
42) 2,4,6-Tribromophenol	10.757	330	26458	135.86	ng	0.00
45) 2-Fluorobiphenyl	9.286	172	144607	84.26	ng	0.00
79) Terphenyl-d14	13.039	244	140771	84.32	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.963	88	8933	41.76	ng	# 100
3) Pyridine	3.675	79	20742	38.63	ng	97
4) n-Nitrosodimethylamine	3.646	42	20895	47.88	ng	# 100
6) Aniline	6.587	93	30150	36.75	ng	96
8) 2-Chlorophenol	6.704	128	35335	49.24	ng	95
9) Benzaldehyde	6.463	77	3349	7.33	ng	92
10) Phenol	6.563	94	38293	50.37	ng	94
11) bis(2-Chloroethyl)ether	6.657	93	29926	50.78	ng	98
12) 1,3-Dichlorobenzene	6.857	146	36942	47.04	ng	96
13) 1,4-Dichlorobenzene	6.934	146	37699	47.00	ng	98
14) 1,2-Dichlorobenzene	7.087	146	36249	47.55	ng	99
15) Benzyl Alcohol	7.063	79	25241	46.85	ng	91
16) 2,2'-oxybis(1-Chloropr...	7.192	45	51217	50.42	ng	98
17) 2-Methylphenol	7.169	107	26575	50.60	ng	98
18) Hexachloroethane	7.434	117	14417	47.73	ng	# 86
19) n-Nitroso-di-n-propyla...	7.340	70	21324	48.71	ng	97
20) 3+4-Methylphenols	7.322	107	33984	49.15	ng	91
22) Acetophenone	7.334	105	48044	47.42	ng	96
24) Nitrobenzene	7.504	77	31246	44.56	ng	99
25) Isophorone	7.745	82	57307	44.14	ng	95
26) 2-Nitrophenol	7.816	139	19660	51.97	ng	# 92
27) 2,4-Dimethylphenol	7.851	122	33235	53.02	ng	95
28) bis(2-Chloroethoxy)met...	7.951	93	37789	49.90	ng	97
29) 2,4-Dichlorophenol	8.057	162	30958	47.88	ng	99
30) 1,2,4-Trichlorobenzene	8.140	180	32733	45.75	ng	97
31) Naphthalene	8.228	128	103902	45.00	ng	97
32) Benzoic acid	8.004	122	23806m	49.78	ng	
33) 4-Chloroaniline	8.269	127	18136	20.79	ng	97
34) Hexachlorobutadiene	8.339	225	18744	46.41	ng	96
35) Caprolactam	8.675	113	9031m	55.22	ng	
36) 4-Chloro-3-methylphenol	8.757	107	30020	46.96	ng	94
37) 2-Methylnaphthalene	8.916	142	71221	45.71	ng	100
38) 1-Methylnaphthalene	9.016	142	65802	44.97	ng	97
40) 1,2,4,5-Tetrachloroben...	9.081	216	31720	47.52	ng	96
41) Hexachlorocyclopentadiene	9.069	237	40245	96.53	ng	97
43) 2,4,6-Trichlorophenol	9.192	196	22861	47.44	ng	98

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.245	196	22109	44.49	ng	98
46) 1,1'-Biphenyl	9.386	154	85760	45.12	ng	99
47) 2-Chloronaphthalene	9.410	162	65720	44.10	ng	99
48) 2-Nitroaniline	9.504	65	17698	49.39	ng	95
49) Acenaphthylene	9.828	152	103766	44.62	ng	99
50) Dimethylphthalate	9.692	163	80375	46.16	ng	97
51) 2,6-Dinitrotoluene	9.751	165	17518	51.04	ng	97
52) Acenaphthene	9.998	154	65191	43.84	ng	99
53) 3-Nitroaniline	9.916	138	10962	28.62	ng	91
54) 2,4-Dinitrophenol	10.028	184	19161	106.98	ng	90
55) Dibenzofuran	10.175	168	95490	43.68	ng	96
56) 4-Nitrophenol	10.086	139	29637	92.97	ng	95
57) 2,4-Dinitrotoluene	10.157	165	23286	49.56	ng	# 92
58) Fluorene	10.516	166	73002	43.17	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.286	232	17806	46.39	ng	# 96
60) Diethylphthalate	10.392	149	83448	45.84	ng	98
61) 4-Chlorophenyl-phenyle...	10.504	204	34394	44.81	ng	96
62) 4-Nitroaniline	10.545	138	17305	44.83	ng	85
63) Azobenzene	10.663	77	68502	43.05	ng	99
65) 4,6-Dinitro-2-methylph...	10.563	198	11673	47.16	ng	96
66) n-Nitrosodiphenylamine	10.628	169	63943	45.32	ng	96
67) 4-Bromophenyl-phenylether	10.992	248	20102	44.87	ng	# 79
68) Hexachlorobenzene	11.057	284	21822	47.54	ng	98
69) Atrazine	11.157	200	19704	45.60	ng	94
70) Pentachlorophenol	11.251	266	27810	95.78	ng	95
71) Phenanthrene	11.480	178	105024	44.43	ng	100
72) Anthracene	11.533	178	107522	45.34	ng	99
73) Carbazole	11.680	167	93854	42.83	ng	99
74) Di-n-butylphthalate	12.010	149	133367	45.18	ng	100
75) Fluoranthene	12.669	202	104886	43.98	ng	97
77) Benzidine	12.780	184	27630	26.36	ng	97
78) Pyrene	12.898	202	104197	44.73	ng	100
80) Butylbenzylphthalate	13.510	149	50919	45.21	ng	98
81) Benzo(a)anthracene	14.080	228	80473	43.28	ng	98
82) 3,3'-Dichlorobenzidine	14.045	252	19806	40.67	ng	97
83) Chrysene	14.121	228	79569	44.67	ng	99
84) Bis(2-ethylhexyl)phtha...	14.063	149	71275	47.14	ng	99
85) Di-n-octyl phthalate	14.680	149	108359	49.53	ng	99
87) Indeno(1,2,3-cd)pyrene	17.098	276	61767	54.02	ng	98
88) Benzo(b)fluoranthene	15.145	252	67067	47.39	ng	98
89) Benzo(k)fluoranthene	15.180	252	64371	49.57	ng	# 96
90) Benzo(a)pyrene	15.521	252	59091	50.58	ng	97
91) Dibenzo(a,h)anthracene	17.121	278	51635	53.05	ng	97
92) Benzo(g,h,i)perylene	17.562	276	50676	53.33	ng	93

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

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