

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF022221\
 Data File : BF123255.D
 Acq On : 22 Feb 2021 12:36
 Operator : JU/CG
 Sample : PB134747BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB134747BS

Manual Integrations
 APPROVED

mohammad
 2/23/2021 3:44:00 PM

Quant Time: Feb 22 13:50:00 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF021921.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 19 16:08:53 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.857	152	114516	20.00	ng	0.00
21) Naphthalene-d8	8.145	136	466159	20.00	ng	0.00
39) Acenaphthene-d10	9.904	164	244460	20.00	ng	0.00
64) Phenanthrene-d10	11.392	188	462269	20.00	ng	0.00
76) Chrysene-d12	14.045	240	389817	20.00	ng	0.00
86) Perylene-d12	15.515	264	355406	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.492	112	1315237	165.32	ng	0.02
7) Phenol-d6	6.498	99	1336389	123.64	ng	0.00
23) Nitrobenzene-d5	7.422	82	894630	87.85	ng	0.00
42) 2,4,6-Tribromophenol	10.698	330	336740	135.28	ng	0.00
45) 2-Fluorobiphenyl	9.222	172	1441022	84.67	ng	0.00
79) Terphenyl-d14	12.986	244	1622968	80.06	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.687	88	180334	39.28	ng	89
3) Pyridine	3.434	79	504613	45.11	ng	93
4) n-Nitrosodimethylamine	3.387	42	266473	51.51	ng	88
6) Aniline	6.522	93	364614	28.56	ng	94
8) 2-Chlorophenol	6.645	128	405636	47.46	ng	93
9) Benzaldehyde	6.404	77	191435	32.71	ng	91
10) Phenol	6.510	94	521842	44.30	ng	94
11) bis(2-Chloroethyl)ether	6.592	93	404462	47.87	ng	96
12) 1,3-Dichlorobenzene	6.798	146	414227	46.59	ng	# 92
13) 1,4-Dichlorobenzene	6.875	146	420691	46.43	ng	94
14) 1,2-Dichlorobenzene	7.028	146	402577	47.76	ng	95
15) Benzyl Alcohol	7.004	79	391984	46.51	ng	93
16) 2,2'-oxybis(1-Chloropr...	7.133	45	521074	44.96	ng	98
17) 2-Methylphenol	7.116	107	339352	47.20	ng	94
18) Hexachloroethane	7.369	117	163450	49.94	ng	93
19) n-Nitroso-di-n-propyla...	7.275	70	308194	47.93	ng	95
20) 3+4-Methylphenols	7.269	107	425714	45.09	ng	# 89
22) Acetophenone	7.269	105	574573	47.52	ng	# 91
24) Nitrobenzene	7.439	77	485362	45.21	ng	90
25) Isophorone	7.681	82	835839	43.89	ng	96
26) 2-Nitrophenol	7.757	139	210123	46.72	ng	# 86
27) 2,4-Dimethylphenol	7.798	122	342054	49.13	ng	93
28) bis(2-Chloroethoxy)met...	7.892	93	496627	49.83	ng	98
29) 2,4-Dichlorophenol	8.004	162	327402	45.55	ng	96
30) 1,2,4-Trichlorobenzene	8.080	180	350328	44.82	ng	99
31) Naphthalene	8.163	128	1140647	44.65	ng	100
32) Benzoic acid	7.933	122	246514	48.00	ng	89
33) 4-Chloroaniline	8.210	127	184457	17.74	ng	# 93
34) Hexachlorobutadiene	8.280	225	194247	44.06	ng	99
35) Caprolactam	8.586	113	114166m	47.33	ng	
36) 4-Chloro-3-methylphenol	8.704	107	406650	47.59	ng	90
37) 2-Methylnaphthalene	8.857	142	776320	45.62	ng	100
38) 1-Methylnaphthalene	8.957	142	722563	45.40	ng	99
40) 1,2,4,5-Tetrachloroben...	9.027	216	333900	45.71	ng	# 97
41) Hexachlorocyclopentadiene	9.016	237	360134	105.23	ng	96
43) 2,4,6-Trichlorophenol	9.139	196	233921	45.15	ng	94

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.180	196	248942	44.90	ng	# 93
46) 1,1'-Biphenyl	9.322	154	947011	46.70	ng	99
47) 2-Chloronaphthalene	9.351	162	714350	44.64	ng	99
48) 2-Nitroaniline	9.445	65	278221	48.53	ng	# 80
49) Acenaphthylene	9.769	152	1099891	45.32	ng	99
50) Dimethylphthalate	9.627	163	857096	46.88	ng	100
51) 2,6-Dinitrotoluene	9.686	165	196763	46.70	ng	# 76
52) Acenaphthene	9.939	154	850195m	50.80	ng	
53) 3-Nitroaniline	9.857	138	118135	25.26	ng	88
54) 2,4-Dinitrophenol	9.963	184	225408	101.85	ng	# 81
55) Dibenzofuran	10.110	168	1005126	44.37	ng	95
56) 4-Nitrophenol	10.027	139	354807	95.21	ng	# 65
57) 2,4-Dinitrotoluene	10.092	165	274515	48.45	ng	# 85
58) Fluorene	10.457	166	815927	45.76	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.227	232	213047	48.17	ng	# 85
60) Diethylphthalate	10.327	149	857691	47.38	ng	99
61) 4-Chlorophenyl-phenyle...	10.445	204	386916	45.58	ng	97
62) 4-Nitroaniline	10.474	138	215257	45.73	ng	84
63) Azobenzene	10.604	77	928328	44.51	ng	95
65) 4,6-Dinitro-2-methylph...	10.504	198	152160	48.90	ng	91
66) n-Nitrosodiphenylamine	10.563	169	722325	45.18	ng	99
67) 4-Bromophenyl-phenylether	10.933	248	229923	45.62	ng	# 89
68) Hexachlorobenzene	11.004	284	247238	46.57	ng	# 90
69) Atrazine	11.092	200	213051	42.17	ng	97
70) Pentachlorophenol	11.204	266	312536	93.24	ng	98
71) Phenanthrene	11.421	178	1247542	45.64	ng	99
72) Anthracene	11.474	178	1293977	47.44	ng	99
73) Carbazole	11.627	167	1154361	44.94	ng	99
74) Di-n-butylphthalate	11.957	149	1345841	46.86	ng	99
75) Fluoranthene	12.610	202	1376436	47.50	ng	100
77) Benzidine	12.733	184	424882	33.74	ng	100
78) Pyrene	12.845	202	1385979	45.54	ng	99
80) Butylbenzylphthalate	13.462	149	599599	46.61	ng	97
81) Benzo(a)anthracene	14.033	228	1220059	46.05	ng	97
82) 3,3'-Dichlorobenzidine	13.998	252	264334	30.02	ng	99
83) Chrysene	14.074	228	1189217	45.71	ng	100
84) Bis(2-ethylhexyl)phtha...	14.021	149	805628	45.49	ng	100
85) Di-n-octyl phthalate	14.639	149	1368612	45.23	ng	97
87) Indeno(1,2,3-cd)pyrene	17.003	276	1138788	46.36	ng	98
88) Benzo(b)fluoranthene	15.092	252	1189396	47.65	ng	99
89) Benzo(k)fluoranthene	15.121	252	1151285	50.87	ng	99
90) Benzo(a)pyrene	15.462	252	1043430	46.88	ng	100
91) Dibenzo(a,h)anthracene	17.027	278	940841	44.65	ng	98
92) Benzo(g,h,i)perylene	17.456	276	913637	47.07	ng	98

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

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