

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF072721\
 Data File : BF124791.D
 Acq On : 27 Jul 2021 13:01
 Operator : JU/CG
 Sample : PB137987BS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB137987BS

Manual Integrations
 APPROVED

mohammad
 7/28/2021 11:32:27 AM

Quant Time: Jul 27 14:43:52 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	8008	20.000	ng	0.00
21) Naphthalene-d8	8.157	136	32844	20.000	ng	0.00
39) Acenaphthene-d10	9.916	164	18570	20.000	ng	0.00
64) Phenanthrene-d10	11.404	188	33381	20.000	ng	0.00
76) Chrysene-d12	14.045	240	23253	20.000	ng	# 0.00
86) Perylene-d12	15.515	264	17230	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.510	112	61572	130.184	ng	0.03
7) Phenol-d6	6.510	99	76312	128.703	ng	0.01
23) Nitrobenzene-d5	7.439	82	41970	76.055	ng	0.00
42) 2,4,6-Tribromophenol	10.710	330	19738	123.795	ng	0.00
45) 2-Fluorobiphenyl	9.233	172	89816	71.038	ng	0.00
79) Terphenyl-d14	12.986	244	92073	67.874	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.810	88	8310	50.265	ng	# 100
3) Pyridine	3.546	79	17681	45.178	ng	97
4) n-Nitrosodimethylamine	3.510	42	15107	48.811	ng	# 89
6) Aniline	6.539	93	24088	39.530	ng	# 94
8) 2-Chlorophenol	6.663	128	25808	48.379	ng	94
9) Benzaldehyde	6.428	77	12417	38.896	ng	93
10) Phenol	6.522	94	30088	52.990	ng	93
11) bis(2-Chloroethyl)ether	6.616	93	24022	53.358	ng	98
12) 1,3-Dichlorobenzene	6.816	146	27737	47.840	ng	98
13) 1,4-Dichlorobenzene	6.892	146	28316	48.818	ng	96
14) 1,2-Dichlorobenzene	7.045	146	26853	47.915	ng	96
15) Benzyl Alcohol	7.016	79	20260	51.961	ng	93
16) 2,2'-oxybis(1-Chloropr...	7.145	45	38084	50.260	ng	96
17) 2-Methylphenol	7.128	107	19795	50.936	ng	97
18) Hexachloroethane	7.386	117	11721	53.314	ng	90
19) n-Nitroso-di-n-propyla...	7.292	70	16785	53.014	ng	91
20) 3+4-Methylphenols	7.281	107	24933	48.559	ng	88
22) Acetophenone	7.286	105	37072	48.708	ng	# 93
24) Nitrobenzene	7.457	77	26013	48.081	ng	90
25) Isophorone	7.698	82	45828	46.966	ng	94
26) 2-Nitrophenol	7.775	139	14678	48.983	ng	94
27) 2,4-Dimethylphenol	7.810	122	24133	52.339	ng	94
28) bis(2-Chloroethoxy)met...	7.904	93	30002	52.910	ng	99
29) 2,4-Dichlorophenol	8.016	162	22725	46.536	ng	95
30) 1,2,4-Trichlorobenzene	8.098	180	24678	46.109	ng	97
31) Naphthalene	8.181	128	77354	45.130	ng	97
32) Benzoic acid	7.951	122	15149m	42.447	ng	
33) 4-Chloroaniline	8.228	127	14499	22.495	ng	98
34) Hexachlorobutadiene	8.292	225	14370	44.916	ng	96
35) Caprolactam	8.616	113	6866m	54.096	ng	
36) 4-Chloro-3-methylphenol	8.710	107	24036	51.604	ng	90
37) 2-Methylnaphthalene	8.875	142	51943	45.355	ng	97
38) 1-Methylnaphthalene	8.975	142	48855	45.030	ng	100
40) 1,2,4,5-Tetrachloroben...	9.039	216	24098	46.510	ng	# 96
41) Hexachlorocyclopentadiene	9.022	237	33511	109.319	ng	98
43) 2,4,6-Trichlorophenol	9.151	196	16899	45.944	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.192	196	17001	45.016	ng	91
46) 1,1'-Biphenyl	9.339	154	63180	45.579	ng	99
47) 2-Chloronaphthalene	9.363	162	48679	44.362	ng	96
48) 2-Nitroaniline	9.457	65	14688	51.119	ng	91
49) Acenaphthylene	9.780	152	77099	45.561	ng	100
50) Dimethylphthalate	9.645	163	60055	46.967	ng	# 99
51) 2,6-Dinitrotoluene	9.704	165	13135	46.498	ng	97
52) Acenaphthene	9.951	154	45699	40.738	ng	97
53) 3-Nitroaniline	9.869	138	8785	29.560	ng	86
54) 2,4-Dinitrophenol	9.980	184	13850	84.788	ng	94
55) Dibenzofuran	10.122	168	68928	43.000	ng	98
56) 4-Nitrophenol	10.039	139	20484	87.282	ng	88
57) 2,4-Dinitrotoluene	10.110	165	18082	47.169	ng	# 94
58) Fluorene	10.469	166	54770	44.539	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.239	232	13628	45.691	ng	# 93
60) Diethylphthalate	10.339	149	62791	47.224	ng	99
61) 4-Chlorophenyl-phenyle...	10.457	204	26376	44.370	ng	97
62) 4-Nitroaniline	10.492	138	12941	44.007	ng	98
63) Azobenzene	10.616	77	55105	46.845	ng	97
65) 4,6-Dinitro-2-methylph...	10.516	198	8499	39.315	ng	98
66) n-Nitrosodiphenylamine	10.580	169	48047	44.416	ng	97
67) 4-Bromophenyl-phenylether	10.945	248	16306	46.059	ng	# 81
68) Hexachlorobenzene	11.016	284	17552	47.738	ng	# 90
69) Atrazine	11.104	200	15866	48.241	ng	98
70) Pentachlorophenol	11.210	266	19154	83.777	ng	93
71) Phenanthrene	11.433	178	79693	44.388	ng	99
72) Anthracene	11.486	178	83550	46.549	ng	99
73) Carbazole	11.633	167	70346	41.812	ng	99
74) Di-n-butylphthalate	11.963	149	105405	47.036	ng	98
75) Fluoranthene	12.621	202	80344	43.853	ng	98
77) Benzidine	12.739	184	37105	56.587	ng	98
78) Pyrene	12.851	202	80755	43.868	ng	98
80) Butylbenzylphthalate	13.463	149	40189	45.581	ng	98
81) Benzo(a)anthracene	14.033	228	64749	42.599	ng	99
82) 3,3'-Dichlorobenzidine	13.998	252	13540	32.098	ng	# 94
83) Chrysene	14.074	228	62785	42.875	ng	97
84) Bis(2-ethylhexyl)phtha...	14.015	149	61334	47.226	ng	99
85) Di-n-octyl phthalate	14.627	149	97067	46.058	ng	98
87) Indeno(1,2,3-cd)pyrene	17.003	276	42173	42.396	ng	95
88) Benzo(b)fluoranthene	15.086	252	50279	41.446	ng	97
89) Benzo(k)fluoranthene	15.121	252	52684	47.530	ng	98
90) Benzo(a)pyrene	15.457	252	47623	46.994	ng	# 93
91) Dibenzo(a,h)anthracene	17.021	278	36172	41.538	ng	98
92) Benzo(g,h,i)perylene	17.451	276	34011	41.187	ng	# 88

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

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