

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM081721\
 Data File : BM031778.D
 Acq On : 17 Aug 2021 12:59
 Operator : CG/JU
 Sample : PB138394BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB138394BS

Manual Integrations
 APPROVED

mohammad
 8/19/2021 1:24:26 PM

Quant Time: Aug 17 14:48:24 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM081621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 16 16:42:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.551	152	45260	20.000	ng	0.00
21) Naphthalene-d8	10.310	136	193955	20.000	ng	# 0.00
39) Acenaphthene-d10	14.186	164	122644	20.000	ng	0.00
64) Phenanthrene-d10	16.939	188	221999	20.000	ng	0.00
76) Chrysene-d12	21.139	240	180835	20.000	ng	0.00
86) Perylene-d12	23.285	264	187293	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.187	112	268217	95.114	ng	0.00
7) Phenol-d6	6.745	99	327509	80.151	ng	0.00
23) Nitrobenzene-d5	8.704	82	386613	83.462	ng	0.00
42) 2,4,6-Tribromophenol	15.686	330	106501	70.506	ng	0.00
45) 2-Fluorobiphenyl	12.798	172	811968	88.120	ng	0.00
79) Terphenyl-d14	19.586	244	989896	92.230	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.193	88	40769	34.580	ng	96
3) Pyridine	3.569	79	97257	29.851	ng	# 81
4) n-Nitrosodimethylamine	3.493	42	82833	43.565	ng	86
6) Aniline	6.898	93	187352	41.933	ng	98
8) 2-Chlorophenol	7.122	128	141300	42.792	ng	95
9) Benzaldehyde	6.710	77	93474	32.230	ng	96
10) Phenol	6.775	94	173722	42.796	ng	89
11) bis(2-Chloroethyl)ether	6.998	93	120795	39.614	ng	94
12) 1,3-Dichlorobenzene	7.439	146	143876	40.735	ng	# 94
13) 1,4-Dichlorobenzene	7.587	146	147993	40.667	ng	95
14) 1,2-Dichlorobenzene	7.892	146	144842	40.519	ng	96
15) Benzyl Alcohol	7.798	79	136580	38.815	ng	96
16) 2,2'-oxybis(1-Chloropr...	8.081	45	165721	38.035	ng	98
17) 2-Methylphenol	7.998	107	118259	41.199	ng	# 89
18) Hexachloroethane	8.604	117	62514	41.441	ng	90
19) n-Nitroso-di-n-propyla...	8.363	70	114601	35.518	ng	94
20) 3+4-Methylphenols	8.328	107	160763	39.843	ng	95
22) Acetophenone	8.369	105	224380	40.742	ng	# 91
24) Nitrobenzene	8.745	77	183377	38.680	ng	94
25) Isophorone	9.269	82	291137	35.041	ng	96
26) 2-Nitrophenol	9.439	139	77025	41.274	ng	# 81
27) 2,4-Dimethylphenol	9.510	122	130064	45.578	ng	91
28) bis(2-Chloroethoxy)met...	9.745	93	170040	42.534	ng	99
29) 2,4-Dichlorophenol	9.969	162	136001	41.946	ng	93
30) 1,2,4-Trichlorobenzene	10.175	180	145225	41.946	ng	96
31) Naphthalene	10.363	128	448336	40.374	ng	99
32) Benzoic acid	9.681	122	77481	30.325	ng	92
33) 4-Chloroaniline	10.486	127	151031	32.721	ng	94
34) Hexachlorobutadiene	10.633	225	91880	41.563	ng	95
35) Caprolactam	11.292	113	35226m	31.290	ng	
36) 4-Chloro-3-methylphenol	11.616	107	148396	36.911	ng	89
37) 2-Methylnaphthalene	11.980	142	324392	39.037	ng	# 96
38) 1-Methylnaphthalene	12.198	142	302505	38.115	ng	# 99
40) 1,2,4,5-Tetrachloroben...	12.351	216	159664	45.832	ng	# 95
41) Hexachlorocyclopentadiene	12.327	237	224847	129.012	ng	99
43) 2,4,6-Trichlorophenol	12.604	196	106160	41.340	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.674	196	114369	40.753	ng	# 88
46) 1,1'-Biphenyl	13.010	154	400594	41.506	ng	99
47) 2-Chloronaphthalene	13.051	162	321006	42.366	ng	# 92
48) 2-Nitroaniline	13.274	65	108685	38.651	ng	89
49) Acenaphthylene	13.904	152	506672	40.845	ng	99
50) Dimethylphthalate	13.663	163	379284	37.114	ng	99
51) 2,6-Dinitrotoluene	13.780	165	82801	36.485	ng	# 70
52) Acenaphthene	14.251	154	296032	39.174	ng	98
53) 3-Nitroaniline	14.110	138	67083	28.788	ng	91
54) 2,4-Dinitrophenol	14.327	184	66184	49.469	ng	# 75
55) Dibenzofuran	14.592	168	479438	37.967	ng	92
56) 4-Nitrophenol	14.433	139	136910	68.877	ng	# 73
57) 2,4-Dinitrotoluene	14.574	165	112341	34.371	ng	# 61
58) Fluorene	15.245	166	394474	37.630	ng	100
59) 2,3,4,6-Tetrachlorophenol	14.821	232	90491	35.559	ng	# 83
60) Diethylphthalate	15.033	149	399081	35.528	ng	97
61) 4-Chlorophenyl-phenyle...	15.245	204	195051	38.546	ng	# 85
62) 4-Nitroaniline	15.286	138	76096	31.783	ng	# 79
63) Azobenzene	15.539	77	452099	36.359	ng	92
65) 4,6-Dinitro-2-methylph...	15.345	198	50787	33.464	ng	78
66) n-Nitrosodiphenylamine	15.462	169	323065	44.264	ng	99
67) 4-Bromophenyl-phenylether	16.139	248	107537	45.508	ng	# 91
68) Hexachlorobenzene	16.245	284	114050	44.122	ng	# 89
69) Atrazine	16.427	200	108476	41.763	ng	99
70) Pentachlorophenol	16.592	266	144601	83.284	ng	96
71) Phenanthrene	16.980	178	553114	40.852	ng	99
72) Anthracene	17.074	178	567623	42.764	ng	99
73) Carbazole	17.351	167	485432	36.506	ng	# 97
74) Di-n-butylphthalate	17.921	149	617026	37.255	ng	# 98
75) Fluoranthene	19.003	202	569208	35.099	ng	98
77) Benzidine	19.209	184	406019	75.554	ng	100
78) Pyrene	19.368	202	575982	43.967	ng	99
80) Butylbenzylphthalate	20.292	149	252140	40.158	ng	94
81) Benzo(a)anthracene	21.127	228	531594	40.133	ng	99
82) 3,3'-Dichlorobenzidine	21.068	252	165900	40.530	ng	99
83) Chrysene	21.174	228	521861	40.419	ng	99
84) Bis(2-ethylhexyl)phtha...	21.068	149	376636	38.818	ng	# 94
85) Di-n-octyl phthalate	21.933	149	626105	39.054	ng	# 94
87) Indeno(1,2,3-cd)pyrene	25.421	276	641995	48.641	ng	99
88) Benzo(b)fluoranthene	22.650	252	567170	40.333	ng	99
89) Benzo(k)fluoranthene	22.697	252	530997	39.122	ng	99
90) Benzo(a)pyrene	23.197	252	543102	43.621	ng	98
91) Dibenzo(a,h)anthracene	25.427	278	548949	47.291	ng	99
92) Benzo(g,h,i)perylene	26.068	276	536517	50.756	ng	98

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

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