

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060721\
 Data File : BP005742.D
 Acq On : 07 Jun 2021 14:56
 Operator : CG/JU
 Sample : M2580-18MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 18-B1(0-2)MSD

Manual Integrations
 APPROVED

mohammad
 6/8/2021 3:37:17 PM

Quant Time: Jun 07 16:31:51 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP052721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 07 12:13:53 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	156967	20.000	ng	0.00
21) Naphthalene-d8	10.787	136	611599	20.000	ng	0.00
39) Acenaphthene-d10	14.604	164	390199	20.000	ng	0.00
64) Phenanthrene-d10	17.345	188	819920	20.000	ng	# 0.00
76) Chrysene-d12	21.416	240	888517	20.000	ng	# 0.00
86) Perylene-d12	23.845	264	1020622	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.546	112	1212190	120.065	ng	0.01
7) Phenol-d6	7.152	99	1412732	102.644	ng	0.01
23) Nitrobenzene-d5	9.146	82	1136483	103.342	ng	0.00
42) 2,4,6-Tribromophenol	16.093	330	877654	175.408	ng	0.00
45) 2-Fluorobiphenyl	13.234	172	2511900	85.195	ng	0.00
79) Terphenyl-d14	19.886	244	3993475	84.337	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.417	88	159304	34.476	ng	# 40
3) Pyridine	3.834	79	310566	27.151	ng	94
4) n-Nitrosodimethylamine	3.728	42	208096	40.847	ng	# 25
6) Aniline	7.305	93	294605	18.311	ng	94
8) 2-Chlorophenol	7.546	128	544042	48.125	ng	97
9) Benzaldehyde	7.117	77	239331	40.163	ng	93
10) Phenol	7.175	94	546630	38.899	ng	97
11) bis(2-Chloroethyl)ether	7.399	93	486672	44.856	ng	97
12) 1,3-Dichlorobenzene	7.869	146	547803	42.690	ng	97
13) 1,4-Dichlorobenzene	8.016	146	562264	43.235	ng	98
14) 1,2-Dichlorobenzene	8.334	146	541525	43.539	ng	99
15) Benzyl Alcohol	8.222	79	490697	49.702	ng	# 83
16) 2,2'-oxybis(1-Chloropr...	8.516	45	495960	39.932	ng	86
17) 2-Methylphenol	8.428	107	423696	46.316	ng	97
18) Hexachloroethane	9.058	117	202852	43.955	ng	92
19) n-Nitroso-di-n-propyla...	8.793	70	386114	44.103	ng	98
20) 3+4-Methylphenols	8.758	107	570499	46.812	ng	95
22) Acetophenone	8.805	105	790332	48.446	ng	# 98
24) Nitrobenzene	9.187	77	584544	47.853	ng	91
25) Isophorone	9.716	82	1042284	42.689	ng	98
26) 2-Nitrophenol	9.899	139	294769	61.805	ng	# 81
27) 2,4-Dimethylphenol	9.958	122	496873	53.020	ng	98
28) bis(2-Chloroethoxy)met...	10.193	93	628433	47.825	ng	97
29) 2,4-Dichlorophenol	10.434	162	526492	52.235	ng	99
30) 1,2,4-Trichlorobenzene	10.652	180	541583	45.365	ng	99
31) Naphthalene	10.834	128	1568904	43.077	ng	98
32) Benzoic acid	10.140	122	292708	55.762	ng	96
33) 4-Chloroaniline	10.952	127	140572	9.269	ng	# 90
34) Hexachlorobutadiene	11.116	225	346526	46.473	ng	98
35) Caprolactam	11.752	113	128320m	46.489	ng	
36) 4-Chloro-3-methylphenol	12.075	107	536334	50.552	ng	93
37) 2-Methylnaphthalene	12.440	142	1155806	44.787	ng	# 88
38) 1-Methylnaphthalene	12.657	142	1075654	43.932	ng	# 95
40) 1,2,4,5-Tetrachloroben...	12.810	216	635532	48.825	ng	96
41) Hexachlorocyclopentadiene	12.787	237	699153	96.480	ng	98
43) 2,4,6-Trichlorophenol	13.046	196	436966	55.348	ng	94

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.122	196	475271	55.669	ng	# 93
46) 1,1'-Biphenyl	13.440	154	1479343	46.691	ng	99
47) 2-Chloronaphthalene	13.487	162	1143964	44.419	ng	99
48) 2-Nitroaniline	13.687	65	344582	55.581	ng	# 76
49) Acenaphthylene	14.328	152	1836674	45.753	ng	99
50) Dimethylphthalate	14.063	163	1625371	52.369	ng	99
51) 2,6-Dinitrotoluene	14.181	165	330721	48.220	ng	# 66
52) Acenaphthene	14.669	154	1082818	41.495	ng	97
53) 3-Nitroaniline	14.510	138	224461	34.238	ng	81
54) 2,4-Dinitrophenol	14.722	184	399131	135.555	ng	98
55) Dibenzofuran	14.998	168	1729250	43.234	ng	99
56) 4-Nitrophenol	14.828	139	513646	88.315	ng	# 60
57) 2,4-Dinitrotoluene	14.969	165	461859	53.890	ng	# 65
58) Fluorene	15.645	166	1389649	43.455	ng	99
59) 2,3,4,6-Tetrachlorophenol	15.228	232	420602m	56.941	ng	
60) Diethylphthalate	15.416	149	1503484	48.281	ng	98
61) 4-Chlorophenyl-phenyle...	15.634	204	720672	45.075	ng	95
62) 4-Nitroaniline	15.669	138	345254	48.536	ng	# 51
63) Azobenzene	15.934	77	1370772	42.036	ng	95
65) 4,6-Dinitro-2-methylph...	15.734	198	261469	57.851	ng	90
66) n-Nitrosodiphenylamine	15.851	169	1248618	44.692	ng	# 92
67) 4-Bromophenyl-phenylether	16.534	248	488657	49.472	ng	98
68) Hexachlorobenzene	16.651	284	578367	49.177	ng	# 93
69) Atrazine	16.804	200	489553	61.314	ng	97
70) Pentachlorophenol	16.998	266	676314	114.658	ng	96
71) Phenanthrene	17.387	178	2252497	44.096	ng	99
72) Anthracene	17.481	178	2308479	45.770	ng	99
73) Carbazole	17.745	167	2130702	42.924	ng	99
74) Di-n-butylphthalate	18.287	149	2598644	48.886	ng	# 92
75) Fluoranthene	19.345	202	2759540	45.004	ng	97
77) Benzidine	19.516	184	150042	8.892	ng	99
78) Pyrene	19.698	202	2847464	45.418	ng	99
80) Butylbenzylphthalate	20.557	149	1228324	48.399	ng	# 90
81) Benzo(a)anthracene	21.398	228	2864083	43.923	ng	98
82) 3,3'-Dichlorobenzidine	21.322	252	646411	30.691	ng	# 97
83) Chrysene	21.451	228	2787731	44.150	ng	99
84) Bis(2-ethylhexyl)phtha...	21.310	149	1757504	49.623	ng	# 96
85) Di-n-octyl phthalate	22.239	149	3210095	53.929	ng	97
87) Indeno(1,2,3-cd)pyrene	26.421	276	3636009	41.873	ng	# 93
88) Benzo(b)fluoranthene	23.104	252	3218385	43.461	ng	# 98
89) Benzo(k)fluoranthene	23.151	252	3052003	43.494	ng	# 96
90) Benzo(a)pyrene	23.739	252	3053453	45.325	ng	# 96
91) Dibenzo(a,h)anthracene	26.439	278	3097848	41.109	ng	# 95
92) Benzo(g,h,i)perylene	27.215	276	2996490	41.226	ng	# 94

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

