

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082321\
 Data File : BP006832.D
 Acq On : 23 Aug 2021 15:46
 Operator : CG/JU
 Sample : M3463-09MS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SWCS-20-0306MS

Quant Time: Aug 23 17:01:50 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080321.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 23 11:39:38 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.693	152	152131	20.000	ng	0.00
21) Naphthalene-d8	10.481	136	651593	20.000	ng	# 0.00
39) Acenaphthene-d10	14.340	164	373763	20.000	ng	0.00
64) Phenanthrene-d10	17.104	188	733146	20.000	ng	# 0.00
76) Chrysene-d12	21.216	240	717401	20.000	ng	# 0.00
86) Perylene-d12	23.527	264	763373	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.299	112	851142	85.933	ng	0.00
7) Phenol-d6	6.887	99	1081369	76.857	ng	0.00
23) Nitrobenzene-d5	8.857	82	727572	51.540	ng	0.00
42) 2,4,6-Tribromophenol	15.845	330	309142	79.141	ng	0.00
45) 2-Fluorobiphenyl	12.963	172	988555	39.571	ng	0.00
79) Terphenyl-d14	19.680	244	994317	27.772	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.240	88	166827	38.688	ng	97
3) Pyridine	3.634	79	409089	32.278	ng	95
4) n-Nitrosodimethylamine	3.546	42	172872	36.138	ng	# 75
6) Aniline	7.034	93	607100	39.723	ng	98
8) 2-Chlorophenol	7.263	128	520232	48.544	ng	96
9) Benzaldehyde	6.846	77	331663	39.165	ng	97
10) Phenol	6.916	94	646124	45.798	ng	97
11) bis(2-Chloroethyl)ether	7.134	93	467142	43.607	ng	97
12) 1,3-Dichlorobenzene	7.581	146	524987	47.003	ng	99
13) 1,4-Dichlorobenzene	7.728	146	533074	47.026	ng	97
14) 1,2-Dichlorobenzene	8.040	146	512134	47.083	ng	95
15) Benzyl Alcohol	7.946	79	482411	45.721	ng	94
16) 2,2'-oxybis(1-Chloropr...	8.222	45	476055	37.344	ng	95
17) 2-Methylphenol	8.152	107	428840	48.138	ng	99
18) Hexachloroethane	8.758	117	203053	45.256	ng	90
19) n-Nitroso-di-n-propyla...	8.510	70	391189	41.442	ng	94
20) 3+4-Methylphenols	8.481	107	586174	48.342	ng	97
22) Acetophenone	8.522	105	774182	44.803	ng	# 98
24) Nitrobenzene	8.899	77	591252	40.291	ng	94
25) Isophorone	9.428	82	1042637	41.121	ng	95
26) 2-Nitrophenol	9.604	139	265704	49.753	ng	94
27) 2,4-Dimethylphenol	9.675	122	476005	53.227	ng	96
28) bis(2-Chloroethoxy)met...	9.910	93	628284	46.311	ng	98
29) 2,4-Dichlorophenol	10.140	162	452129	49.937	ng	95
30) 1,2,4-Trichlorobenzene	10.346	180	450530	46.308	ng	98
31) Naphthalene	10.534	128	1547303	44.192	ng	99
32) Benzoic acid	9.857	122	385568	55.732	ng	94
33) 4-Chloroaniline	10.652	127	330679	23.339	ng	96
34) Hexachlorobutadiene	10.810	225	269050	47.457	ng	99
35) Caprolactam	11.475	113	164525	50.395	ng	# 81
36) 4-Chloro-3-methylphenol	11.799	107	550729	47.875	ng	91
37) 2-Methylnaphthalene	12.151	142	1083441	45.665	ng	99
38) 1-Methylnaphthalene	12.369	142	999707	44.339	ng	98
40) 1,2,4,5-Tetrachloroben...	12.522	216	476168	48.676	ng	# 96
41) Hexachlorocyclopentadiene	12.493	237	421482	81.236	ng	99
43) 2,4,6-Trichlorophenol	12.775	196	339338	50.802	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	12.846	196	367999	49.772	ng	# 91
46) 1,1'-Biphenyl	13.175	154	1274016	45.034	ng	98
47) 2-Chloronaphthalene	13.210	162	1005456	46.152	ng	94
48) 2-Nitroaniline	13.434	65	359090	46.705	ng	92
49) Acenaphthylene	14.063	152	1638050	46.630	ng	100
50) Dimethylphthalate	13.816	163	1262691	46.411	ng	99
51) 2,6-Dinitrotoluene	13.934	165	272610	44.904	ng	# 81
52) Acenaphthene	14.404	154	955757	44.705	ng	99
53) 3-Nitroaniline	14.263	138	232712	34.577	ng	84
54) 2,4-Dinitrophenol	14.481	184	239179	75.337	ng	# 85
55) Dibenzofuran	14.745	168	1492123	44.371	ng	94
56) 4-Nitrophenol	14.592	139	552275	94.502	ng	87
57) 2,4-Dinitrotoluene	14.728	165	380762	46.766	ng	# 79
58) Fluorene	15.398	166	1219140	45.357	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.975	232	301064	48.589	ng	# 74
60) Diethylphthalate	15.187	149	1339464	46.874	ng	97
61) 4-Chlorophenyl-phenyle...	15.398	204	565728	46.476	ng	91
62) 4-Nitroaniline	15.434	138	276492	38.715	ng	93
63) Azobenzene	15.692	77	1416355	42.512	ng	93
65) 4,6-Dinitro-2-methylph...	15.492	198	150208	33.531	ng	77
66) n-Nitrosodiphenylamine	15.616	169	1035718	44.945	ng	99
67) 4-Bromophenyl-phenylether	16.292	248	339507	48.263	ng	# 87
68) Hexachlorobenzene	16.404	284	379166	47.365	ng	# 91
69) Atrazine	16.586	200	376216	52.319	ng	96
70) Pentachlorophenol	16.757	266	514764	101.350	ng	100
71) Phenanthrene	17.145	178	1862177	45.030	ng	99
72) Anthracene	17.239	178	1902077	47.600	ng	100
73) Carbazole	17.516	167	1665920	40.873	ng	98
74) Di-n-butylphthalate	18.081	149	2320016	49.961	ng	99
75) Fluoranthene	19.128	202	2115377	45.575	ng	95
77) Benzidine	19.322	184	438191	24.409	ng	100
78) Pyrene	19.480	202	2176565	45.422	ng	99
80) Butylbenzylphthalate	20.369	149	1123310	53.310	ng	90
81) Benzo(a)anthracene	21.198	228	2088160	44.538	ng	100
82) 3,3'-Dichlorobenzidine	21.139	252	759244	61.684	ng	# 98
83) Chrysene	21.251	228	2030398	44.671	ng	99
84) Bis(2-ethylhexyl)phtha...	21.133	149	1644808	51.944	ng	# 97
85) Di-n-octyl phthalate	22.033	149	2911275	55.781	ng	98
87) Indeno(1,2,3-cd)pyrene	25.933	276	2585671	46.713	ng	# 92
88) Benzo(b)fluoranthene	22.827	252	2183421	44.009	ng	# 97
89) Benzo(k)fluoranthene	22.874	252	2132229	44.762	ng	# 97
90) Benzo(a)pyrene	23.433	252	2136870	48.019	ng	# 96
91) Dibenzo(a,h)anthracene	25.951	278	2201516	46.001	ng	# 94
92) Benzo(g,h,i)perylene	26.668	276	2159596	47.089	ng	# 92

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

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