

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071824\
 Data File : BM046713.D
 Acq On : 18 Jul 2024 22:01
 Operator : MA/JU
 Sample : P3246-06MS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 WC-CTW2-BL-02MS

Quant Time: Jul 19 01:16:15 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM071724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 18 06:01:46 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.487	152	116468	20.000	ng	0.00
21) Naphthalene-d8	10.246	136	423278	20.000	ng	0.00
39) Acenaphthene-d10	14.134	164	278352	20.000	ng	0.00
64) Phenanthrene-d10	16.886	188	511900	20.000	ng	0.00
76) Chrysene-d12	21.116	240	370439	20.000	ng	0.00
86) Perylene-d12	23.874	264	472284	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.134	112	673311	117.423	ng	0.00
7) Phenol-d6	6.693	99	810679	111.028	ng	0.00
23) Nitrobenzene-d5	8.634	82	581039	71.271	ng	0.00
42) 2,4,6-Tribromophenol	15.634	330	415067	96.136	ng	0.00
45) 2-Fluorobiphenyl	12.746	172	1516464	76.955	ng	0.00
79) Terphenyl-d14	19.539	244	1640576	80.089	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.140	88	69003	39.651	ng	94
3) Pyridine	3.511	79	194930	38.219	ng	96
4) n-Nitrosodimethylamine	3.429	42	192449	50.940	ng	99
6) Aniline	6.828	93	250994	39.774	ng	99
8) 2-Chlorophenol	7.064	128	342645	53.385	ng	99
9) Benzaldehyde	6.640	77	28334	6.357	ng	94
10) Phenol	6.717	94	359182	50.619	ng	95
11) bis(2-Chloroethyl)ether	6.928	93	271271	51.458	ng	100
12) 1,3-Dichlorobenzene	7.375	146	413002	48.720	ng	99
13) 1,4-Dichlorobenzene	7.522	146	416951	48.430	ng	98
14) 1,2-Dichlorobenzene	7.828	146	398532	48.592	ng	100
15) Benzyl Alcohol	7.734	79	310578	52.225	ng	100
16) 2,2'-oxybis(1-Chloropr...	8.011	45	300660	54.557	ng	97
17) 2-Methylphenol	7.940	107	253243	50.107	ng	97
18) Hexachloroethane	8.546	117	160096	51.207	ng	97
19) n-Nitroso-di-n-propyla...	8.299	70	231382	50.199	ng	98
20) 3+4-Methylphenols	8.269	107	343881	49.524	ng	98
22) Acetophenone	8.305	105	450546	46.290	ng	98
24) Nitrobenzene	8.675	77	380066	47.543	ng	97
25) Isophorone	9.199	82	612951	49.991	ng	99
26) 2-Nitrophenol	9.375	139	196871	51.413	ng	96
27) 2,4-Dimethylphenol	9.458	122	228621	54.988	ng	98
28) bis(2-Chloroethoxy)met...	9.687	93	358074	51.228	ng	99
29) 2,4-Dichlorophenol	9.910	162	373633	50.032	ng	98
30) 1,2,4-Trichlorobenzene	10.116	180	448036	48.974	ng	100
31) Naphthalene	10.299	128	1029411	48.956	ng	99
32) Benzoic acid	9.616	122	189700	37.460	ng	96
33) 4-Chloroaniline	10.416	127	135752	17.327	ng	97
34) Hexachlorobutadiene	10.587	225	299570	48.406	ng	98
35) Caprolactam	11.216	113	72756	36.787	ng	98
36) 4-Chloro-3-methylphenol	11.563	107	314327	45.620	ng	100
37) 2-Methylnaphthalene	11.922	142	763326	47.805	ng	99
38) 1-Methylnaphthalene	12.146	142	705025	44.947	ng	100
40) 1,2,4,5-Tetrachloroben...	12.299	216	465212	51.395	ng	# 99
41) Hexachlorocyclopentadiene	12.281	237	567541	256.268	ng	95
43) 2,4,6-Trichlorophenol	12.551	196	318639	53.479	ng	97

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44) 2,4,5-Trichlorophenol	12.628	196	328651	49.471	ng	98
46) 1,1'-Biphenyl	12.957	154	964623	49.811	ng	99
47) 2-Chloronaphthalene	12.993	162	835254	52.166	ng	99
48) 2-Nitroaniline	13.210	65	207091	49.295	ng	98
49) Acenaphthylene	13.851	152	1244469	55.945	ng	99
50) Dimethylphthalate	13.610	163	966894	50.228	ng	98
51) 2,6-Dinitrotoluene	13.722	165	207226	46.568	ng	99
52) Acenaphthene	14.198	154	721329	50.567	ng	99
53) 3-Nitroaniline	14.045	138	133882	33.242	ng	98
54) 2,4-Dinitrophenol	14.257	184	248842	94.821	ng	98
55) Dibenzofuran	14.540	168	1201237	49.754	ng	98
56) 4-Nitrophenol	14.387	139	289782	82.911	ng	93
57) 2,4-Dinitrotoluene	14.516	165	280428	43.515	ng	# 98
58) Fluorene	15.187	166	981030	47.543	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.769	232	284636	46.521	ng	100
60) Diethylphthalate	14.987	149	928423	46.752	ng	99
61) 4-Chlorophenyl-phenyle...	15.192	204	530011	47.759	ng	98
62) 4-Nitroaniline	15.216	138	177796	39.153	ng	98
63) Azobenzene	15.487	77	786658	48.789	ng	98
65) 4,6-Dinitro-2-methylph...	15.281	198	171253	64.381	ng	97
66) n-Nitrosodiphenylamine	15.410	169	812290	61.188	ng	97
67) 4-Bromophenyl-phenylether	16.087	248	332988	57.599	ng	97
68) Hexachlorobenzene	16.192	284	377333	54.041	ng	99
69) Atrazine	16.381	200	254785	54.680	ng	95
70) Pentachlorophenol	16.545	266	457937	92.443	ng	99
71) Phenanthrene	16.928	178	1357980	53.419	ng	99
72) Anthracene	17.022	178	1353613	53.819	ng	99
73) Carbazole	17.292	167	1131530	47.543	ng	99
74) Di-n-butylphthalate	17.886	149	1306979	49.117	ng	99
75) Fluoranthene	18.957	202	1373680	42.051	ng	100
77) Benzidine	19.151	184	429299	57.915	ng	99
78) Pyrene	19.322	202	1371244	51.997	ng	99
80) Butylbenzylphthalate	20.251	149	474489	49.811	ng	99
81) Benzo(a)anthracene	21.098	228	1297533	52.931	ng	100
82) 3,3'-Dichlorobenzidine	21.033	252	407775	47.866	ng	99
83) Chrysene	21.157	228	1211732	53.017	ng	100
84) Bis(2-ethylhexyl)phtha...	21.057	149	630256	49.611	ng	99
85) Di-n-octyl phthalate	22.116	149	1066327	49.798	ng	99
87) Indeno(1,2,3-cd)pyrene	26.956	276	1853611	60.578	ng	# 97
88) Benzo(b)fluoranthene	23.015	252	1480917	50.928	ng	100
89) Benzo(k)fluoranthene	23.074	252	1391586	50.702	ng	99
90) Benzo(a)pyrene	23.751	252	1403127	55.778	ng	98
91) Dibenzo(a,h)anthracene	27.015	278	1486235	58.520	ng	100
92) Benzo(g,h,i)perylene	27.915	276	1475326	56.295	ng	98

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

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