

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082321\
 Data File : BF125171.D
 Acq On : 23 Aug 2021 23:59
 Operator : JU/CG
 Sample : M3419-09MS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SWCS-03-1015MS

Manual Integrations
 APPROVED

mohammad
 8/24/2021 11:20:08 AM

Quant Time: Aug 24 04:50:06 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF081821.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 23 10:15:49 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.922	152	164905	20.000	ng	0.00
21) Naphthalene-d8	8.210	136	644596	20.000	ng	# 0.00
39) Acenaphthene-d10	9.963	164	321817	20.000	ng	0.00
64) Phenanthrene-d10	11.451	188	511585	20.000	ng	0.00
76) Chrysene-d12	14.098	240	347816	20.000	ng	# 0.00
86) Perylene-d12	15.592	264	273521	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.551	112	1147529	105.327	ng	0.01
7) Phenol-d6	6.551	99	1467011	104.113	ng	0.00
23) Nitrobenzene-d5	7.486	82	1019061	79.590	ng	0.00
42) 2,4,6-Tribromophenol	10.757	330	395258	123.040	ng	0.00
45) 2-Fluorobiphenyl	9.286	172	1636113	70.857	ng	0.00
79) Terphenyl-d14	13.033	244	1387078	63.532	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.793	88	206119	38.340	ng	# 100
3) Pyridine	3.551	79	472275	33.277	ng	88
4) n-Nitrosodimethylamine	3.510	42	324277	45.662	ng	# 51
6) Aniline	6.587	93	656405	39.659	ng	# 93
8) 2-Chlorophenol	6.710	128	501784	45.148	ng	85
9) Benzaldehyde	6.475	77	384320	38.865	ng	93
10) Phenol	6.563	94	663926	44.994	ng	91
11) bis(2-Chloroethyl)ether	6.657	93	511109	43.980	ng	85
12) 1,3-Dichlorobenzene	6.863	146	555960	43.167	ng	97
13) 1,4-Dichlorobenzene	6.939	146	563334	43.897	ng	97
14) 1,2-Dichlorobenzene	7.092	146	541729	44.870	ng	98
15) Benzyl Alcohol	7.063	79	497038	45.814	ng	96
16) 2,2'-oxybis(1-Chloropr...	7.198	45	1016428	43.612	ng	99
17) 2-Methylphenol	7.175	107	414369	44.382	ng	95
18) Hexachloroethane	7.439	117	205635	45.760	ng	# 84
19) n-Nitroso-di-n-propyla...	7.339	70	373297	44.045	ng	# 80
20) 3+4-Methylphenols	7.322	107	491740	41.959	ng	92
22) Acetophenone	7.334	105	699089	42.141	ng	93
24) Nitrobenzene	7.510	77	594663	42.624	ng	90
25) Isophorone	7.745	82	1024979	41.497	ng	# 89
26) 2-Nitrophenol	7.822	139	271642	48.761	ng	# 81
27) 2,4-Dimethylphenol	7.857	122	461498	47.320	ng	86
28) bis(2-Chloroethoxy)met...	7.951	93	620694	45.385	ng	# 97
29) 2,4-Dichlorophenol	8.063	162	431492	43.383	ng	96
30) 1,2,4-Trichlorobenzene	8.145	180	463673	42.751	ng	96
31) Naphthalene	8.228	128	1347487	40.818	ng	99
32) Benzoic acid	7.981	122	342465	51.736	ng	# 81
33) 4-Chloroaniline	8.275	127	397061	30.510	ng	# 92
34) Hexachlorobutadiene	8.345	225	297362	42.997	ng	99
35) Caprolactam	8.651	113	140441m	54.515	ng	
36) 4-Chloro-3-methylphenol	8.757	107	465108	45.806	ng	94
37) 2-Methylnaphthalene	8.922	142	938149	43.006	ng	98
38) 1-Methylnaphthalene	9.022	142	859461	42.191	ng	96
40) 1,2,4,5-Tetrachloroben...	9.086	216	467624	44.475	ng	98
41) Hexachlorocyclopentadiene	9.069	237	336939	60.995	ng	98
43) 2,4,6-Trichlorophenol	9.198	196	321530	43.431	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.233	196	344116	44.181	ng	98
46) 1,1'-Biphenyl	9.386	154	1069599	41.225	ng	99
47) 2-Chloronaphthalene	9.410	162	886714	42.056	ng	99
48) 2-Nitroaniline	9.504	65	325429	49.793	ng	# 79
49) Acenaphthylene	9.827	152	1304303	42.453	ng	98
50) Dimethylphthalate	9.686	163	1043443	46.788	ng	# 98
51) 2,6-Dinitrotoluene	9.745	165	225427	46.616	ng	# 77
52) Acenaphthene	9.998	154	866139	45.476	ng	98
53) 3-Nitroaniline	9.916	138	183603	35.207	ng	# 78
54) 2,4-Dinitrophenol	10.022	184	168404	88.163	ng	# 86
55) Dibenzofuran	10.175	168	1112233	40.577	ng	99
56) 4-Nitrophenol	10.075	139	343354	83.143	ng	# 82
57) 2,4-Dinitrotoluene	10.151	165	296334	50.632	ng	# 95
58) Fluorene	10.516	166	849241	41.907	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.286	232	254090	43.959	ng	# 87
60) Diethylphthalate	10.380	149	1012012	46.547	ng	99
61) 4-Chlorophenyl-phenyle...	10.504	204	443231	43.259	ng	89
62) 4-Nitroaniline	10.533	138	198564	42.343	ng	86
63) Azobenzene	10.663	77	977851	40.098	ng	92
65) 4,6-Dinitro-2-methylph...	10.557	198	114601	42.325	ng	86
66) n-Nitrosodiphenylamine	10.622	169	826574	45.123	ng	95
67) 4-Bromophenyl-phenylether	10.992	248	289053	47.052	ng	97
68) Hexachlorobenzene	11.063	284	297066	47.006	ng	# 82
69) Atrazine	11.151	200	257442	48.977	ng	90
70) Pentachlorophenol	11.251	266	319750	86.651	ng	93
71) Phenanthrene	11.480	178	1181906	41.929	ng	97
72) Anthracene	11.527	178	1195465	43.027	ng	98
73) Carbazole	11.680	167	1036095	40.807	ng	100
74) Di-n-butylphthalate	12.010	149	1337660	44.397	ng	100
75) Fluoranthene	12.668	202	1186487	42.420	ng	97
77) Benzidine	12.786	184	674340	55.513	ng	99
78) Pyrene	12.898	202	1190641	39.907	ng	96
80) Butylbenzylphthalate	13.510	149	523798	44.679	ng	92
81) Benzo(a)anthracene	14.086	228	1060462	42.316	ng	98
82) 3,3'-Dichlorobenzidine	14.045	252	390964	50.329	ng	98
83) Chrysene	14.121	228	1035321	41.775	ng	97
84) Bis(2-ethylhexyl)phtha...	14.068	149	716698	44.480	ng	# 99
85) Di-n-octyl phthalate	14.686	149	1183944	44.179	ng	99
87) Indeno(1,2,3-cd)pyrene	17.115	276	746448	37.881	ng	# 89
88) Benzo(b)fluoranthene	15.151	252	954810	47.781	ng	# 95
89) Benzo(k)fluoranthene	15.186	252	844740	45.302	ng	98
90) Benzo(a)pyrene	15.533	252	821624	46.547	ng	# 95
91) Dibenzo(a,h)anthracene	17.133	278	624176	36.645	ng	# 94
92) Benzo(g,h,i)perylene	17.580	276	612169	37.274	ng	# 87

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

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