

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF012822\
 Data File : BF126719.D
 Acq On : 28 Jan 2022 16:38
 Operator : CG/JU
 Sample : N1319-04MSD
 Misc :
 ALS Vial : 6 Sample Multiplier: 2

Instrument :
 BNA_F
 ClientSampleId :
 MH-1DMSD

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 01/31/2022
 Supervised By : Jagrut Upadhyay 01/31/2022

Quant Time: Jan 29 02:27:09 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF011822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 27 03:09:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.957	152	263494	40.000	ng	0.00
21) Naphthalene-d8	8.240	136	1017729	40.000	ng	0.00
39) Acenaphthene-d10	9.998	164	511234	40.000	ng	0.00
64) Phenanthrene-d10	11.492	188	736272	40.000	ng	0.00
76) Chrysene-d12	14.145	240	483591	40.000	ng	# 0.00
86) Perylene-d12	15.663	264	566968	40.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.587	112	1270619	126.911	ng	0.02
7) Phenol-d6	6.581	99	1729337	124.320	ng	0.00
23) Nitrobenzene-d5	7.522	82	1227951	94.452	ng	0.00
42) 2,4,6-Tribromophenol	10.792	330	289793	121.142	ng	0.00
45) 2-Fluorobiphenyl	9.316	172	1490026	89.273	ng	0.00
79) Terphenyl-d14	13.080	244	1194466	82.857	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	2.869	88	270337	54.786	ng	# 87
3) Pyridine	3.628	79	667869	48.317	ng	# 82
4) n-Nitrosodimethylamine	3.593	42	246812	50.464	ng	# 70
6) Aniline	6.622	93	733926m	42.104	ng	
8) 2-Chlorophenol	6.740	128	524401	57.105	ng	95
9) Benzaldehyde	6.510	77	521594	57.154	ng	86
10) Phenol	6.593	94	843160	56.990	ng	99
11) bis(2-Chloroethyl)ether	6.693	93	689774	57.453	ng	93
12) 1,3-Dichlorobenzene	6.898	146	559452	54.878	ng	96
13) 1,4-Dichlorobenzene	6.975	146	561996	54.591	ng	95
14) 1,2-Dichlorobenzene	7.128	146	545657	56.223	ng	99
15) Benzyl Alcohol	7.093	79	598370	48.249	ng	92
16) 2,2'-oxybis(1-Chloropr...	7.228	45	782627	57.137	ng	85
17) 2-Methylphenol	7.204	107	494706	54.795	ng	97
18) Hexachloroethane	7.469	117	227529	54.738	ng	93
19) n-Nitroso-di-n-propyla...	7.369	70	483942	47.667	ng	# 84
20) 3+4-Methylphenols	7.351	107	598047	51.087	ng	# 81
22) Acetophenone	7.363	105	807249	51.477	ng	# 90
24) Nitrobenzene	7.540	77	769457	57.332	ng	# 88
25) Isophorone	7.775	82	1353138	53.669	ng	95
26) 2-Nitrophenol	7.851	139	262553	74.993	ng	# 76
27) 2,4-Dimethylphenol	7.887	122	494164	59.478	ng	91
28) bis(2-Chloroethoxy)met...	7.987	93	825456	52.494	ng	97
29) 2,4-Dichlorophenol	8.092	162	413335	52.446	ng	97
30) 1,2,4-Trichlorobenzene	8.181	180	446289	52.805	ng	98
31) Naphthalene	8.263	128	1474467	53.637	ng	99
32) Benzoic acid	7.992	122	289805	61.490	ng	# 70
33) 4-Chloroaniline	8.304	127	171847	15.647	ng	# 92
34) Hexachlorobutadiene	8.375	225	270246	50.268	ng	98
35) Caprolactam	8.681	113	138238m	49.118	ng	
36) 4-Chloro-3-methylphenol	8.781	107	497871	48.308	ng	88
37) 2-Methylnaphthalene	8.951	142	914962	53.519	ng	97
38) 1-Methylnaphthalene	9.051	142	848239	51.190	ng	99
40) 1,2,4,5-Tetrachloroben...	9.116	216	419688	58.537	ng	97
41) Hexachlorocyclopentadiene	9.104	237	525916	120.479	ng	95
43) 2,4,6-Trichlorophenol	9.228	196	275007	53.806	ng	96

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.269	196	280769	53.041	ng	# 91
46) 1,1'-Biphenyl	9.422	154	1078959	56.027	ng	98
47) 2-Chloronaphthalene	9.445	162	839448	55.483	ng	100
48) 2-Nitroaniline	9.539	65	328521	62.887	ng	92
49) Acenaphthylene	9.863	152	1323862	55.780	ng	99
50) Dimethylphthalate	9.716	163	924435	50.990	ng	99
51) 2,6-Dinitrotoluene	9.781	165	203862	65.073	ng	90
52) Acenaphthene	10.034	154	851598	58.692	ng	96
53) 3-Nitroaniline	9.951	138	119955	33.490	ng	89
54) 2,4-Dinitrophenol	10.051	184	176429	128.047	ng	95
55) Dibenzofuran	10.210	168	1130564	51.849	ng	98
56) 4-Nitrophenol	10.104	139	294313	103.188	ng	# 78
57) 2,4-Dinitrotoluene	10.186	165	255170	67.216	ng	# 93
58) Fluorene	10.551	166	828341	50.316	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.322	232	220478	50.945	ng	# 82
60) Diethylphthalate	10.416	149	849363	47.129	ng	98
61) 4-Chlorophenyl-phenyle...	10.539	204	404774	49.230	ng	89
62) 4-Nitroaniline	10.569	138	183517	52.792	ng	81
63) Azobenzene	10.704	77	1241254	46.540	ng	92
65) 4,6-Dinitro-2-methylph...	10.586	198	115514	70.335	ng	88
66) n-Nitrosodiphenylamine	10.657	169	706515	56.855	ng	99
67) 4-Bromophenyl-phenylether	11.033	248	245803	57.934	ng	94
68) Hexachlorobenzene	11.098	284	267124	59.010	ng	# 86
69) Atrazine	11.186	200	211077	53.185	ng	92
70) Pentachlorophenol	11.286	266	264723	100.568	ng	97
71) Phenanthrene	11.516	178	1139747	54.191	ng	99
72) Anthracene	11.569	178	1165261	55.220	ng	99
73) Carbazole	11.722	167	951995	48.052	ng	99
74) Di-n-butylphthalate	12.045	149	1173907	49.137	ng	98
75) Fluoranthene	12.710	202	998780	48.078	ng	99
77) Benzidine	12.827	184	487493	58.443	ng	99
78) Pyrene	12.939	202	1004292	51.361	ng	99
80) Butylbenzylphthalate	13.557	149	426607	50.218	ng	# 98
81) Benzo(a)anthracene	14.133	228	893998	54.397	ng	100
82) 3,3'-Dichlorobenzidine	14.092	252	186206	34.489	ng	# 96
83) Chrysene	14.169	228	840375	53.142	ng	99
84) Bis(2-ethylhexyl)phtha...	14.110	149	609227	52.954	ng	99
85) Di-n-octyl phthalate	14.733	149	1089211	61.836	ng	95
87) Indeno(1,2,3-cd)pyrene	17.227	276	1121412	64.025	ng	# 92
88) Benzo(b)fluoranthene	15.210	252	992200	52.779	ng	# 94
89) Benzo(k)fluoranthene	15.245	252	890668	48.254	ng	# 95
90) Benzo(a)pyrene	15.598	252	946562	61.884	ng	# 96
91) Dibenzo(a,h)anthracene	17.245	278	949724	65.452	ng	95
92) Benzo(g,h,i)perylene	17.704	276	973618	68.430	ng	# 92

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

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