

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122021\
 Data File : BF126278.D
 Acq On : 20 Dec 2021 15:04
 Operator : JU/CG
 Sample : M5146-01MS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NB-8-121721MS

Quant Time: Dec 21 05:09:28 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF121521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 16 12:22:48 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.816	152	153602	20.000	ng	0.00
21) Naphthalene-d8	8.104	136	585698	20.000	ng	# 0.00
39) Acenaphthene-d10	9.857	164	228964	20.000	ng	0.00
64) Phenanthrene-d10	11.339	188	331540	20.000	ng	# 0.00
76) Chrysene-d12	13.980	240	268713	20.000	ng	0.01
86) Perylene-d12	15.421	264	296192	20.000	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.434	112	571127	54.756	ng	0.01
7) Phenol-d6	6.445	99	682944	51.567	ng	0.00
23) Nitrobenzene-d5	7.375	82	416626	38.525	ng	0.00
42) 2,4,6-Tribromophenol	10.639	330	94083	51.130	ng	0.00
45) 2-Fluorobiphenyl	9.175	172	568247	43.528	ng	0.00
79) Terphenyl-d14	12.921	244	462707	29.929	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.634	88	268359	52.708	ng	# 100
3) Pyridine	3.363	79	577888	42.740	ng	# 78
4) n-Nitrosodimethylamine	3.316	42	280879	54.052	ng	# 1
6) Aniline	6.475	93	635200	37.759	ng	96
8) 2-Chlorophenol	6.598	128	540197	51.072	ng	88
9) Benzaldehyde	6.363	77	415390	51.977	ng	95
10) Phenol	6.457	94	750903	50.510	ng	90
11) bis(2-Chloroethyl)ether	6.551	93	591580	51.266	ng	97
12) 1,3-Dichlorobenzene	6.757	146	539703	50.583	ng	98
13) 1,4-Dichlorobenzene	6.834	146	534409	49.599	ng	94
14) 1,2-Dichlorobenzene	6.987	146	507452	50.937	ng	95
15) Benzyl Alcohol	6.957	79	455360	47.098	ng	94
16) 2,2'-oxybis(1-Chloropr...	7.098	45	932936	49.404	ng	93
17) 2-Methylphenol	7.075	107	432007	46.623	ng	97
18) Hexachloroethane	7.334	117	208113	49.104	ng	97
19) n-Nitroso-di-n-propyla...	7.234	70	354806	43.640	ng	# 72
20) 3+4-Methylphenols	7.222	107	475090	43.370	ng	# 80
22) Acetophenone	7.228	105	685841	50.880	ng	97
24) Nitrobenzene	7.398	77	567621	52.578	ng	91
25) Isophorone	7.639	82	1008991	49.407	ng	# 87
26) 2-Nitrophenol	7.716	139	258145	53.043	ng	# 83
27) 2,4-Dimethylphenol	7.757	122	444307	53.759	ng	98
28) bis(2-Chloroethoxy)met...	7.851	93	653048	48.638	ng	# 98
29) 2,4-Dichlorophenol	7.957	162	349755	47.608	ng	92
30) 1,2,4-Trichlorobenzene	8.045	180	380214	50.034	ng	97
31) Naphthalene	8.122	128	1370803	50.047	ng	98
32) Benzoic acid	7.869	122	288232	40.481	ng	91
33) 4-Chloroaniline	8.169	127	240623	21.040	ng	98
34) Hexachlorobutadiene	8.245	225	184786	49.512	ng	98
35) Caprolactam	8.539	113	119733	65.578	ng	94
36) 4-Chloro-3-methylphenol	8.651	107	392463	45.288	ng	89
37) 2-Methylnaphthalene	8.816	142	813913	48.252	ng	98
38) 1-Methylnaphthalene	8.916	142	752144	45.947	ng	97
40) 1,2,4,5-Tetrachloroben...	8.980	216	274039	55.454	ng	# 96
41) Hexachlorocyclopentadiene	8.969	237	229178	80.987	ng	94
43) 2,4,6-Trichlorophenol	9.092	196	194567	50.051	ng	95

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF122021\
 Data File : BF126278.D
 Acq On : 20 Dec 2021 15:04
 Operator : JU/CG
 Sample : M5146-01MS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NB-8-121721MS

Quant Time: Dec 21 05:09:28 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF121521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 16 12:22:48 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.128	196	196881	48.907	ng	94
46) 1,1'-Biphenyl	9.280	154	921694	54.442	ng	98
47) 2-Chloronaphthalene	9.304	162	664029	52.131	ng	99
48) 2-Nitroaniline	9.392	65	240924	51.820	ng	89
49) Acenaphthylene	9.716	152	1009837	51.883	ng	98
50) Dimethylphthalate	9.580	163	739426	50.982	ng	99
51) 2,6-Dinitrotoluene	9.639	165	159578	50.833	ng	96
52) Acenaphthene	9.892	154	644486	51.059	ng	99
53) 3-Nitroaniline	9.804	138	137416	34.440	ng #	81
54) 2,4-Dinitrophenol	9.910	184	70947	54.887	ng #	76
55) Dibenzofuran	10.063	168	841154	48.947	ng	95
56) 4-Nitrophenol	9.963	139	283860	98.489	ng	78
57) 2,4-Dinitrotoluene	10.039	165	199325	49.775	ng #	95
58) Fluorene	10.404	166	580853	46.688	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.175	232	139033	46.580	ng #	99
60) Diethylphthalate	10.280	149	696337	47.775	ng	97
61) 4-Chlorophenyl-phenyle...	10.398	204	260156	46.443	ng	90
62) 4-Nitroaniline	10.410	138	143339	42.824	ng #	74
63) Azobenzene	10.557	77	825375	46.752	ng	83
65) 4,6-Dinitro-2-methylph...	10.445	198	46932	25.923	ng	87
66) n-Nitrosodiphenylamine	10.510	169	536212	50.395	ng	97
67) 4-Bromophenyl-phenylether	10.886	248	153867	49.450	ng	94
68) Hexachlorobenzene	10.951	284	177948	49.725	ng	90
69) Atrazine	11.039	200	155318	79.654	ng	97
70) Pentachlorophenol	11.145	266	190000	95.701	ng	92
71) Phenanthrene	11.363	178	861393	50.339	ng	98
72) Anthracene	11.416	178	883900	51.463	ng	99
73) Carbazole	11.569	167	808216	49.968	ng	99
74) Di-n-butylphthalate	11.904	149	1032041	50.332	ng	100
75) Fluoranthene	12.551	202	852944	55.324	ng	91
77) Benzidine	12.674	184	764496	180.459	ng	98
78) Pyrene	12.780	202	900869	36.491	ng	97
80) Butylbenzylphthalate	13.404	149	457403	39.373	ng	88
81) Benzo(a)anthracene	13.968	228	720641	45.275	ng	99
82) 3,3'-Dichlorobenzidine	13.927	252	329155	45.265	ng #	97
83) Chrysene	14.004	228	781418	47.250	ng	98
84) Bis(2-ethylhexyl)phtha...	13.963	149	558984	43.059	ng	99
85) Di-n-octyl phthalate	14.574	149	1241395	53.642	ng	98
87) Indeno(1,2,3-cd)pyrene	16.856	276	848271	42.937	ng #	92
88) Benzo(b)fluoranthene	15.004	252	954614	53.640	ng #	96
89) Benzo(k)fluoranthene	15.033	252	827083	48.581	ng #	95
90) Benzo(a)pyrene	15.362	252	868922	58.745	ng #	94
91) Dibenzo(a,h)anthracene	16.880	278	717082	44.650	ng #	96
92) Benzo(g,h,i)perylene	17.292	276	689660	41.505	ng	92

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

