

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012121\
 Data File : BM028481.D
 Acq On : 21 Jan 2021 10:06
 Operator : CG/JU
 Sample : PB134193BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB134193BS

Quant Time: Jan 21 10:40:13 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM012021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 20 14:45:01 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.028	152	43565	20.00	ng	0.00
21) Naphthalene-d8	10.851	136	188002	20.00	ng	# 0.00
39) Acenaphthene-d10	14.675	164	102189	20.00	ng	0.00
64) Phenanthrene-d10	17.398	188	175267	20.00	ng	# 0.00
76) Chrysene-d12	21.527	240	127840	20.00	ng	# 0.00
86) Perylene-d12	23.815	264	115019	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.552	112	360621	130.76	ng	0.00
7) Phenol-d6	7.199	99	462649	123.21	ng	0.01
23) Nitrobenzene-d5	9.210	82	291699	73.59	ng	0.00
42) 2,4,6-Tribromophenol	16.157	330	148344	109.56	ng	0.00
45) 2-Fluorobiphenyl	13.298	172	640844	78.68	ng	0.00
79) Terphenyl-d14	19.992	244	580835	82.05	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.328	88	34283	30.98	ng	# 65
3) Pyridine	3.763	79	103769	33.88	ng	# 49
4) n-Nitrosodimethylamine	3.669	42	69763	39.61	ng	# 58
6) Aniline	7.351	93	96008	23.60	ng	98
8) 2-Chlorophenol	7.593	128	145578	47.47	ng	97
9) Benzaldehyde	7.157	77	17468	8.94	ng	87
10) Phenol	7.222	94	170029	48.05	ng	88
11) bis(2-Chloroethyl)ether	7.451	93	128051	44.87	ng	91
12) 1,3-Dichlorobenzene	7.916	146	151027	41.74	ng	# 93
13) 1,4-Dichlorobenzene	8.069	146	154966	42.35	ng	98
14) 1,2-Dichlorobenzene	8.387	146	148463	42.29	ng	98
15) Benzyl Alcohol	8.281	79	122549	46.84	ng	# 87
16) 2,2'-oxybis(1-Chloropr...	8.569	45	215316	41.80	ng	71
17) 2-Methylphenol	8.481	107	117537	47.96	ng	# 89
18) Hexachloroethane	9.116	117	57935	42.06	ng	91
19) n-Nitroso-di-n-propyla...	8.851	70	103746	46.40	ng	# 66
20) 3+4-Methylphenols	8.816	107	158672	48.28	ng	90
22) Acetophenone	8.869	105	200576	41.15	ng	# 89
24) Nitrobenzene	9.251	77	154998	40.51	ng	# 87
25) Isophorone	9.781	82	270346	44.52	ng	95
26) 2-Nitrophenol	9.963	139	77448	44.73	ng	# 83
27) 2,4-Dimethylphenol	10.022	122	130488	51.82	ng	89
28) bis(2-Chloroethoxy)met...	10.257	93	175516	40.94	ng	99
29) 2,4-Dichlorophenol	10.498	162	131714	43.84	ng	98
30) 1,2,4-Trichlorobenzene	10.710	180	137808	38.62	ng	98
31) Naphthalene	10.898	128	435591	41.01	ng	100
32) Benzoic acid	10.192	122	92239	40.71	ng	# 86
33) 4-Chloroaniline	11.010	127	42619	9.66	ng	# 91
34) Hexachlorobutadiene	11.175	225	81475	38.27	ng	99
35) Caprolactam	11.816	113	36434	37.47	ng	# 60
36) 4-Chloro-3-methylphenol	12.134	107	134382	43.02	ng	94
37) 2-Methylnaphthalene	12.504	142	306997	41.93	ng	96
38) 1-Methylnaphthalene	12.722	142	285422	41.08	ng	100
40) 1,2,4,5-Tetrachloroben...	12.869	216	144951	43.01	ng	# 98
41) Hexachlorocyclopentadiene	12.845	237	194739	123.68	ng	99
43) 2,4,6-Trichlorophenol	13.110	196	99809	44.15	ng	97

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012121\
 Data File : BM028481.D
 Acq On : 21 Jan 2021 10:06
 Operator : CG/JU
 Sample : PB134193BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB134193BS

Quant Time: Jan 21 10:40:13 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM012021.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 20 14:45:01 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	13.186	196	104574	44.70	ng	97
46) 1,1'-Biphenyl	13.510	154	377379	43.27	ng	99
47) 2-Chloronaphthalene	13.557	162	290009	40.69	ng	99
48) 2-Nitroaniline	13.757	65	85161	38.11	ng	91
49) Acenaphthylene	14.398	152	447494	42.31	ng	100
50) Dimethylphthalate	14.133	163	332299	39.56	ng	99
51) 2,6-Dinitrotoluene	14.257	165	73898	41.35	ng	89
52) Acenaphthene	14.739	154	270727	41.23	ng	98
53) 3-Nitroaniline	14.580	138	30581	15.18	ng	89
54) 2,4-Dinitrophenol	14.792	184	88136	82.80	ng	93
55) Dibenzofuran	15.069	168	414584	41.05	ng	100
56) 4-Nitrophenol	14.886	139	118432	74.65	ng	# 78
57) 2,4-Dinitrotoluene	15.039	165	94968	39.89	ng	92
58) Fluorene	15.716	166	336732	40.70	ng	98
59) 2,3,4,6-Tetrachlorophenol	15.292	232	84172	40.70	ng	93
60) Diethylphthalate	15.486	149	330719	38.78	ng	98
61) 4-Chlorophenyl-phenyle...	15.704	204	162483	39.87	ng	93
62) 4-Nitroaniline	15.739	138	67938	30.97	ng	# 82
63) Azobenzene	15.998	77	328897	41.08	ng	90
65) 4,6-Dinitro-2-methylph...	15.798	198	53530	49.16	ng	# 80
66) n-Nitrosodiphenylamine	15.921	169	281626	47.98	ng	98
67) 4-Bromophenyl-phenylether	16.592	248	95919	46.21	ng	94
68) Hexachlorobenzene	16.710	284	106439	44.13	ng	96
69) Atrazine	16.857	200	72395	41.61	ng	97
70) Pentachlorophenol	17.051	266	123398	93.87	ng	99
71) Phenanthrene	17.445	178	456726	42.37	ng	99
72) Anthracene	17.533	178	469591	45.10	ng	98
73) Carbazole	17.798	167	391400	39.64	ng	99
74) Di-n-butylphthalate	18.345	149	497992	41.92	ng	99
75) Fluoranthene	19.433	202	453650	37.47	ng	96
77) Benzidine	19.615	184	90557	31.50	ng	100
78) Pyrene	19.792	202	450754	49.17	ng	99
80) Butylbenzylphthalate	20.668	149	192012	47.51	ng	97
81) Benzo(a)anthracene	21.515	228	374994	42.48	ng	99
82) 3,3'-Dichlorobenzidine	21.445	252	66107	23.30	ng	# 99
83) Chrysene	21.562	228	362739	42.56	ng	100
84) Bis(2-ethylhexyl)phtha...	21.427	149	281522	49.26	ng	97
85) Di-n-octyl phthalate	22.315	149	458117	47.42	ng	# 90
87) Indeno(1,2,3-cd)pyrene	26.150	276	390035	45.51	ng	# 90
88) Benzo(b)fluoranthene	23.127	252	371990	47.64	ng	# 97
89) Benzo(k)fluoranthene	23.174	252	338253	44.36	ng	# 97
90) Benzo(a)pyrene	23.715	252	318350	44.00	ng	# 97
91) Dibenzo(a,h)anthracene	26.162	278	327971	46.26	ng	# 95
92) Benzo(g,h,i)perylene	26.868	276	329366	47.23	ng	# 92

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

