

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060122\
 Data File : BF128619.D
 Acq On : 01 Jun 2022 12:18
 Operator : CG\JU
 Sample : PB145210BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB145210BS

Quant Time: Jun 02 04:34:58 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 27 01:51:59 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.834	152	152965	20.000	ng	0.00
21) Naphthalene-d8	8.122	136	608043	20.000	ng	# 0.00
39) Acenaphthene-d10	9.875	164	351673	20.000	ng	0.00
64) Phenanthrene-d10	11.363	188	639456	20.000	ng	0.00
76) Chrysene-d12	14.010	240	422323	20.000	ng	# 0.00
86) Perylene-d12	15.468	264	289833	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	1067462	115.592	ng	0.05
7) Phenol-d6	6.516	99	1289250	109.494	ng	0.05
23) Nitrobenzene-d5	7.404	82	989072	93.838	ng	0.00
42) 2,4,6-Tribromophenol	10.675	330	456110	136.261	ng	0.00
45) 2-Fluorobiphenyl	9.198	172	1804519	83.364	ng	0.00
79) Terphenyl-d14	12.951	244	2022866	92.538	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.734	88	109306	26.545	ng	96
3) Pyridine	3.469	79	284392	26.676	ng	90
4) n-Nitrosodimethylamine	3.434	42	251120	44.951	ng	# 76
6) Aniline	6.504	93	471173	34.292	ng	# 88
8) 2-Chlorophenol	6.646	128	416030	43.447	ng	99
9) Benzaldehyde	6.387	77	233866	47.711	ng	94
10) Phenol	6.528	94	484464m	41.522	ng	
11) bis(2-Chloroethyl)ether	6.569	93	400036	41.883	ng	99
12) 1,3-Dichlorobenzene	6.775	146	459271	42.215	ng	95
13) 1,4-Dichlorobenzene	6.851	146	457307	41.558	ng	98
14) 1,2-Dichlorobenzene	7.004	146	431811	42.635	ng	97
15) Benzyl Alcohol	6.993	79	393978	43.871	ng	94
16) 2,2'-oxybis(1-Chloropr...	7.110	45	620825	36.727	ng	92
17) 2-Methylphenol	7.134	107	329083	40.379	ng	94
18) Hexachloroethane	7.345	117	182017	46.892	ng	93
19) n-Nitroso-di-n-propyla...	7.263	70	330362	45.760	ng	99
20) 3+4-Methylphenols	7.281	107	468516	47.585	ng	# 80
22) Acetophenone	7.251	105	613460	43.755	ng	# 97
24) Nitrobenzene	7.422	77	480119	46.076	ng	95
25) Isophorone	7.663	82	833655	42.875	ng	96
26) 2-Nitrophenol	7.740	139	214495	49.251	ng	# 92
27) 2,4-Dimethylphenol	7.798	122	362627	41.890	ng	93
28) bis(2-Chloroethoxy)met...	7.869	93	486882	38.583	ng	99
29) 2,4-Dichlorophenol	8.004	162	360295	40.217	ng	99
30) 1,2,4-Trichlorobenzene	8.063	180	389615	40.150	ng	100
31) Naphthalene	8.145	128	1200290	40.333	ng	99
32) Benzoic acid	7.951	122	201643	32.952	ng	92
33) 4-Chloroaniline	8.198	127	371089	31.317	ng	97
34) Hexachlorobutadiene	8.257	225	274914	44.869	ng	99
35) Caprolactam	8.581	113	107199m	38.805	ng	
36) 4-Chloro-3-methylphenol	8.710	107	420358	45.674	ng	92
37) 2-Methylnaphthalene	8.834	142	774717	41.242	ng	100
38) 1-Methylnaphthalene	8.934	142	742103	40.619	ng	98
40) 1,2,4,5-Tetrachloroben...	8.998	216	399793	41.399	ng	99
41) Hexachlorocyclopentadiene	8.987	237	527113	97.102	ng	98
43) 2,4,6-Trichlorophenol	9.128	196	276236	38.789	ng	96

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060122\
 Data File : BF128619.D
 Acq On : 01 Jun 2022 12:18
 Operator : CG\JU
 Sample : PB145210BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB145210BS

Quant Time: Jun 02 04:34:58 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 27 01:51:59 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.187	196	273245m	37.239	ng	
46) 1,1'-Biphenyl	9.298	154	984162	39.855	ng	99
47) 2-Chloronaphthalene	9.322	162	752781	38.834	ng	97
48) 2-Nitroaniline	9.428	65	270568	50.195	ng	94
49) Acenaphthylene	9.739	152	1225443	41.151	ng	100
50) Dimethylphthalate	9.610	163	951069	41.692	ng	99
51) 2,6-Dinitrotoluene	9.669	165	203869	48.764	ng	88
52) Acenaphthene	9.910	154	855783m	46.578	ng	
53) 3-Nitroaniline	9.839	138	168879	34.978	ng	97
54) 2,4-Dinitrophenol	9.945	184	216263	100.112	ng	# 89
55) Dibenzofuran	10.086	168	1057095	38.778	ng	94
56) 4-Nitrophenol	10.051	139	319394	81.522	ng	# 65
57) 2,4-Dinitrotoluene	10.069	165	266254	53.232	ng	# 83
58) Fluorene	10.428	166	850414	43.333	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.210	232	241363	40.590	ng	98
60) Diethylphthalate	10.304	149	946149	42.697	ng	99
61) 4-Chlorophenyl-phenyle...	10.416	204	435189	43.707	ng	91
62) 4-Nitroaniline	10.469	138	198343m	42.907	ng	
63) Azobenzene	10.581	77	890908	39.491	ng	98
65) 4,6-Dinitro-2-methylph...	10.481	198	143083	48.371	ng	99
66) n-Nitrosodiphenylamine	10.545	169	774987	40.480	ng	100
67) 4-Bromophenyl-phenylether	10.910	248	281993	42.683	ng	98
68) Hexachlorobenzene	10.975	284	318102	43.600	ng	# 93
69) Atrazine	11.075	200	275349	46.342	ng	98
70) Pentachlorophenol	11.181	266	311486	70.060	ng	99
71) Phenanthrene	11.392	178	1247808	41.193	ng	99
72) Anthracene	11.445	178	1260894	41.672	ng	100
73) Carbazole	11.604	167	1086285	37.461	ng	99
74) Di-n-butylphthalate	11.928	149	1429532	41.885	ng	99
75) Fluoranthene	12.580	202	1317251	40.674	ng	97
77) Benzidine	12.698	184	451517	46.352	ng	100
78) Pyrene	12.810	202	1295986	40.734	ng	99
80) Butylbenzylphthalate	13.427	149	551852	41.122	ng	99
81) Benzo(a)anthracene	13.998	228	1056735	42.098	ng	98
82) 3,3'-Dichlorobenzidine	13.963	252	302704	47.094	ng	97
83) Chrysene	14.033	228	1002770	39.008	ng	100
84) Bis(2-ethylhexyl)phtha...	13.986	149	762535	46.980	ng	99
85) Di-n-octyl phthalate	14.598	149	1166813	40.839	ng	100
87) Indeno(1,2,3-cd)pyrene	16.933	276	726334	42.489	ng	# 93
88) Benzo(b)fluoranthene	15.045	252	855832	47.468	ng	98
89) Benzo(k)fluoranthene	15.074	252	828727	49.610	ng	97
90) Benzo(a)pyrene	15.410	252	749527	51.815	ng	# 98
91) Dibenzo(a,h)anthracene	16.957	278	642389	45.322	ng	# 95
92) Benzo(g,h,i)perylene	17.386	276	637910	45.375	ng	# 92

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF060122\
 Data File : BF128619.D
 Acq On : 01 Jun 2022 12:18
 Operator : CG\JU
 Sample : PB145210BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB145210BS

Quant Time: Jun 02 04:34:58 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF052622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 27 01:51:59 2022
 Response via : Initial Calibration

