

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101722\
 Data File : BF130685.D
 Acq On : 17 Oct 2022 14:17
 Operator : CG\JU
 Sample : PB148379BSD
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB148379BSD

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 10/18/2022
 Supervised By : Jagrut Upadhyay 10/18/2022

Quant Time: Oct 18 03:20:41 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 18 03:15:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.575	152	81336	20.000	ng	0.00
21) Naphthalene-d8	7.857	136	317683	20.000	ng	0.00
39) Acenaphthene-d10	9.610	164	166169	20.000	ng	0.00
64) Phenanthrene-d10	11.086	188	299023	20.000	ng	0.00
76) Chrysene-d12	13.722	240	233113	20.000	ng	0.00
86) Perylene-d12	15.092	264	157123	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.187	112	644465	142.824	ng	0.02
7) Phenol-d6	6.234	99	775871	136.761	ng	0.00
23) Nitrobenzene-d5	7.151	82	486919	83.672	ng	0.00
42) 2,4,6-Tribromophenol	10.398	330	250856	154.282	ng	0.00
45) 2-Fluorobiphenyl	8.939	172	980899	89.416	ng	0.00
79) Terphenyl-d14	12.675	244	1193011	84.941	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.210	88	71966	26.362	ng	99
3) Pyridine	2.875	79	177712	36.393	ng	98
4) n-Nitrosodimethylamine	2.834	42	109958	47.792	ng	95
6) Aniline	6.240	93	267647	39.406	ng	# 79
8) 2-Chlorophenol	6.363	128	250432	47.194	ng	97
9) Benzaldehyde	6.122	77	139129	39.034	ng	95
10) Phenol	6.245	94	295944	45.084	ng	91
11) bis(2-Chloroethyl)ether	6.322	93	220613	46.267	ng	96
12) 1,3-Dichlorobenzene	6.510	146	266728	45.908	ng	98
13) 1,4-Dichlorobenzene	6.593	146	275980	46.822	ng	97
14) 1,2-Dichlorobenzene	6.740	146	261720	47.677	ng	99
15) Benzyl Alcohol	6.734	79	182593	44.829	ng	98
16) 2,2'-oxybis(1-Chloropr...	6.863	45	267695	43.638	ng	87
17) 2-Methylphenol	6.851	107	190060	45.027	ng	96
18) Hexachloroethane	7.081	117	102749	46.079	ng	99
19) n-Nitroso-di-n-propyla...	7.004	70	160561	43.866	ng	99
20) 3+4-Methylphenols	7.004	107	246053	43.502	ng	97
22) Acetophenone	6.998	105	345302	45.310	ng	# 97
24) Nitrobenzene	7.169	77	249673	42.662	ng	98
25) Isophorone	7.410	82	437862	45.208	ng	98
26) 2-Nitrophenol	7.481	139	133202	46.412	ng	90
27) 2,4-Dimethylphenol	7.528	122	202163	50.540	ng	100
28) bis(2-Chloroethoxy)met...	7.622	93	264785	47.214	ng	99
29) 2,4-Dichlorophenol	7.728	162	202949	45.348	ng	97
30) 1,2,4-Trichlorobenzene	7.804	180	215821	44.732	ng	99
31) Naphthalene	7.881	128	713856	43.410	ng	99
32) Benzoic acid	7.698	122	106272	41.995	ng	95
33) 4-Chloroaniline	7.940	127	212127	33.300	ng	100
34) Hexachlorobutadiene	7.992	225	138777	46.343	ng	99
35) Caprolactam	8.322	113	65349m	48.532	ng	
36) 4-Chloro-3-methylphenol	8.439	107	215022	44.294	ng	99
37) 2-Methylnaphthalene	8.569	142	468221	45.245	ng	99
38) 1-Methylnaphthalene	8.669	142	438755	43.674	ng	99
40) 1,2,4,5-Tetrachloroben...	8.734	216	219120	48.603	ng	99
41) Hexachlorocyclopentadiene	8.722	237	157277	89.871	ng	99
43) 2,4,6-Trichlorophenol	8.857	196	133886	44.229	ng	99

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101722\
 Data File : BF130685.D
 Acq On : 17 Oct 2022 14:17
 Operator : CG\JU
 Sample : PB148379BSD
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB148379BSD

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 10/18/2022
 Supervised By : Jagrut Upadhyay 10/18/2022

Quant Time: Oct 18 03:20:41 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101122.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 18 03:15:54 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	8.904	196	146928	44.136	ng	96
46) 1,1'-Biphenyl	9.034	154	572095	47.506	ng	99
47) 2-Chloronaphthalene	9.057	162	447747	45.492	ng	98
48) 2-Nitroaniline	9.169	65	133231	44.968	ng	99
49) Acenaphthylene	9.469	152	664383	45.052	ng	99
50) Dimethylphthalate	9.345	163	519921	45.200	ng	99
51) 2,6-Dinitrotoluene	9.410	165	116721	46.259	ng	98
52) Acenaphthene	9.639	154	404859	45.313	ng	99
53) 3-Nitroaniline	9.581	138	95516	34.534	ng	99
54) 2,4-Dinitrophenol	9.698	184	114983	88.843	ng	98
55) Dibenzofuran	9.816	168	617398	45.360	ng	99
56) 4-Nitrophenol	9.763	139	134155	87.970	ng	94
57) 2,4-Dinitrotoluene	9.816	165	159391	49.152	ng	91
58) Fluorene	10.157	166	508555	45.447	ng	99
59) 2,3,4,6-Tetrachlorophenol	9.939	232	121172	45.657	ng	96
60) Diethylphthalate	10.045	149	509670	44.409	ng	98
61) 4-Chlorophenyl-phenyle...	10.151	204	234778	46.407	ng	94
62) 4-Nitroaniline	10.198	138	120332	45.568	ng	98
63) Azobenzene	10.310	77	459585	42.516	ng	98
65) 4,6-Dinitro-2-methylph...	10.228	198	84007	44.956	ng	99
66) n-Nitrosodiphenylamine	10.275	169	436639	43.898	ng	99
67) 4-Bromophenyl-phenylether	10.633	248	153601	45.398	ng	95
68) Hexachlorobenzene	10.698	284	179669	46.976	ng	93
69) Atrazine	10.810	200	147185	54.341	ng	97
70) Pentachlorophenol	10.904	266	137467	92.348	ng	97
71) Phenanthrene	11.116	178	742609	44.825	ng	100
72) Anthracene	11.163	178	742690	45.902	ng	99
73) Carbazole	11.328	167	696813	48.264	ng	99
74) Di-n-butylphthalate	11.657	149	807538	46.175	ng	100
75) Fluoranthene	12.298	202	799815	50.232	ng	99
77) Benzidine	12.427	184	314653	53.401	ng	98
78) Pyrene	12.522	202	811327	40.343	ng	99
80) Butylbenzylphthalate	13.151	149	345739	44.709	ng	99
81) Benzo(a)anthracene	13.710	228	691743	44.039	ng	100
82) 3,3'-Dichlorobenzidine	13.680	252	197254	43.458	ng	# 97
83) Chrysene	13.745	228	650023	43.647	ng	99
84) Bis(2-ethylhexyl)phtha...	13.710	149	443246	47.279	ng	99
85) Di-n-octyl phthalate	14.316	149	648844	45.190	ng	99
87) Indeno(1,2,3-cd)pyrene	16.380	276	533711	45.383	ng	98
88) Benzo(b)fluoranthene	14.716	252	542363	54.512	ng	99
89) Benzo(k)fluoranthene	14.739	252	507117	50.058	ng	98
90) Benzo(a)pyrene	15.039	252	426303	51.482	ng	98
91) Dibenzo(a,h)anthracene	16.392	278	465695	48.310	ng	99
92) Benzo(g,h,i)perylene	16.768	276	480505	53.519	ng	98

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101722\
Data File : BF130685.D
Acq On : 17 Oct 2022 14:17
Operator : CG\JU
Sample : PB148379BSD
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PB148379BSD

Manual Integrations APPROVED

Reviewed By :Christian Giraldo 10/18/2022
Supervised By :Jaqrut Upadhyay 10/18/2022

Quant Time: Oct 18 03:20:41 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF101122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 18 03:15:54 2022
Response via : Initial Calibration

