

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020422\
 Data File : BF126805.D
 Acq On : 04 Feb 2022 12:35
 Operator : CG/JU
 Sample : PB142443BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB142443BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 02/07/2022
 Supervised By : Jagrut Upadhyay 02/07/2022

Quant Time: Feb 04 23:52:58 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF020222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 02 15:08:47 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.951	152	157306	20.000	ng	0.00
21) Naphthalene-d8	8.233	136	616825	20.000	ng	0.00
39) Acenaphthene-d10	9.992	164	336061	20.000	ng	0.00
64) Phenanthrene-d10	11.486	188	556458	20.000	ng	0.00
76) Chrysene-d12	14.133	240	382402	20.000	ng	# 0.00
86) Perylene-d12	15.645	264	273463	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.581	112	1434424	124.895	ng	0.02
7) Phenol-d6	6.575	99	1944503	122.397	ng	0.00
23) Nitrobenzene-d5	7.516	82	1389119	84.283	ng	0.00
42) 2,4,6-Tribromophenol	10.786	330	423839	121.522	ng	0.00
45) 2-Fluorobiphenyl	9.310	172	1689738	83.256	ng	0.00
79) Terphenyl-d14	13.074	244	1921808	84.070	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.904	88	218632	37.619	ng	# 86
3) Pyridine	3.651	79	567948	36.036	ng	# 81
4) n-Nitrosodimethylamine	3.616	42	219740	45.138	ng	# 64
6) Aniline	6.616	93	825621	40.727	ng	97
8) 2-Chlorophenol	6.734	128	500691	47.677	ng	97
9) Benzaldehyde	6.504	77	457487	45.772	ng	88
10) Phenol	6.586	94	765723	45.031	ng	97
11) bis(2-Chloroethyl)ether	6.686	93	635652	46.400	ng	91
12) 1,3-Dichlorobenzene	6.892	146	511781	43.280	ng	94
13) 1,4-Dichlorobenzene	6.969	146	518169	43.355	ng	96
14) 1,2-Dichlorobenzene	7.122	146	505537	43.527	ng	100
15) Benzyl Alcohol	7.092	79	563877	44.423	ng	93
16) 2,2'-oxybis(1-Chloropr...	7.222	45	674557	45.706	ng	82
17) 2-Methylphenol	7.192	107	469874	45.499	ng	# 94
18) Hexachloroethane	7.463	117	210269	44.912	ng	# 89
19) n-Nitroso-di-n-propyla...	7.363	70	452548	43.260	ng	# 84
20) 3+4-Methylphenols	7.345	107	558388	42.930	ng	# 84
22) Acetophenone	7.363	105	754601	42.278	ng	# 87
24) Nitrobenzene	7.533	77	718031	43.887	ng	# 87
25) Isophorone	7.769	82	1287644	45.522	ng	# 94
26) 2-Nitrophenol	7.845	139	251270	46.081	ng	# 73
27) 2,4-Dimethylphenol	7.881	122	478339	50.227	ng	88
28) bis(2-Chloroethoxy)met...	7.981	93	771345	42.997	ng	97
29) 2,4-Dichlorophenol	8.086	162	409607	43.690	ng	97
30) 1,2,4-Trichlorobenzene	8.169	180	429023	42.553	ng	97
31) Naphthalene	8.257	128	1383719	42.919	ng	99
32) Benzoic acid	8.004	122	322361	43.861	ng	# 76
33) 4-Chloroaniline	8.298	127	260790	20.020	ng	# 92
34) Hexachlorobutadiene	8.369	225	260758	41.262	ng	99
35) Caprolactam	8.675	113	151340m	44.364	ng	
36) 4-Chloro-3-methylphenol	8.780	107	514817	43.081	ng	89
37) 2-Methylnaphthalene	8.945	142	866080	43.595	ng	97
38) 1-Methylnaphthalene	9.045	142	817030	42.725	ng	99
40) 1,2,4,5-Tetrachloroben...	9.110	216	409073	42.277	ng	96
41) Hexachlorocyclopentadiene	9.098	237	551575	98.897	ng	96
43) 2,4,6-Trichlorophenol	9.222	196	286220	41.989	ng	97

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020422\
 Data File : BF126805.D
 Acq On : 04 Feb 2022 12:35
 Operator : CG/JU
 Sample : PB142443BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB142443BS

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 02/07/2022
 Supervised By : Jagrut Upadhyay 02/07/2022

Quant Time: Feb 04 23:52:58 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF020222.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 02 15:08:47 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.257	196	301278	42.335	ng	94
46) 1,1'-Biphenyl	9.410	154	1048427	42.597	ng	96
47) 2-Chloronaphthalene	9.439	162	822799	42.172	ng	99
48) 2-Nitroaniline	9.533	65	348262	43.780	ng	93
49) Acenaphthylene	9.857	152	1319557	43.804	ng	99
50) Dimethylphthalate	9.716	163	966837	41.621	ng	# 99
51) 2,6-Dinitrotoluene	9.775	165	216078	45.283	ng	86
52) Acenaphthene	10.027	154	890808	46.380	ng	96
53) 3-Nitroaniline	9.945	138	162428	29.872	ng	# 81
54) 2,4-Dinitrophenol	10.051	184	224943	89.492	ng	95
55) Dibenzofuran	10.204	168	1146911	40.456	ng	98
56) 4-Nitrophenol	10.104	139	353252	85.783	ng	# 78
57) 2,4-Dinitrotoluene	10.180	165	288281	44.575	ng	95
58) Fluorene	10.545	166	864689	41.224	ng	99
59) 2,3,4,6-Tetrachlorophenol	10.310	232	255806	41.530	ng	# 81
60) Diethylphthalate	10.416	149	935123	40.970	ng	97
61) 4-Chlorophenyl-phenyle...	10.533	204	426178	40.168	ng	91
62) 4-Nitroaniline	10.569	138	221817	42.990	ng	# 77
63) Azobenzene	10.692	77	1303212	41.151	ng	91
65) 4,6-Dinitro-2-methylph...	10.586	198	143982	42.653	ng	95
66) n-Nitrosodiphenylamine	10.657	169	782484	43.727	ng	99
67) 4-Bromophenyl-phenylether	11.027	248	278031	43.652	ng	# 88
68) Hexachlorobenzene	11.092	284	304794	43.733	ng	# 85
69) Atrazine	11.180	200	267617	46.089	ng	94
70) Pentachlorophenol	11.286	266	348434	85.692	ng	98
71) Phenanthrene	11.510	178	1303258	42.287	ng	99
72) Anthracene	11.563	178	1345902	43.602	ng	100
73) Carbazole	11.716	167	1193386	40.981	ng	99
74) Di-n-butylphthalate	12.039	149	1436971	42.561	ng	# 98
75) Fluoranthene	12.704	202	1344447	42.888	ng	97
77) Benzidine	12.821	184	721740	82.272	ng	99
78) Pyrene	12.933	202	1356798	44.397	ng	99
80) Butylbenzylphthalate	13.545	149	571426	45.064	ng	# 90
81) Benzo(a)anthracene	14.121	228	1117109	43.780	ng	98
82) 3,3'-Dichlorobenzidine	14.080	252	208778	28.570	ng	# 97
83) Chrysene	14.162	228	1063873	44.036	ng	100
84) Bis(2-ethylhexyl)phtha...	14.098	149	754254	45.719	ng	100
85) Di-n-octyl phthalate	14.721	149	1083532	46.062	ng	95
87) Indeno(1,2,3-cd)pyrene	17.198	276	779072	46.709	ng	# 90
88) Benzo(b)fluoranthene	15.198	252	831945	46.527	ng	# 94
89) Benzo(k)fluoranthene	15.227	252	830561	46.565	ng	# 95
90) Benzo(a)pyrene	15.580	252	771653	54.137	ng	# 95
91) Dibenzo(a,h)anthracene	17.221	278	670786	49.099	ng	# 93
92) Benzo(g,h,i)perylene	17.674	276	674529	49.987	ng	# 88

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

```
Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020422\  
Data File : BF126805.D  
Acq On    : 04 Feb 2022  12:35  
Operator  : CG/JU  
Sample    : PB142443BS  
Misc      :  
ALS Vial  : 6    Sample Multiplier: 1
```

Instrument :
BNA_F
ClientSampleId :
PB142443BS

Manual Integrations

Reviewed By :Christian Giraldo	02/07/2022
Supervised By :Jagrut Upadhyay	02/07/2022

Quant Time: Feb 04 23:52:58 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF020222.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 02 15:08:47 2022
Response via : Initial Calibration

