

Data Path : Z:\svoasrv\HPCHEM1\BNA\_F\Data\BF041923\  
 Data File : BF132918.D  
 Acq On : 19 Apr 2023 18:34  
 Operator : CG\JU  
 Sample : 02270-26MS  
 Misc :  
 ALS Vial : 6 Sample Multiplier: 1

Instrument :  
 BNA\_F  
 ClientSampleId :  
 SB-16-14-16-20230413MS

Manual Integrations  
 APPROVED

Reviewed By : Christian Giraldo 04/21/2023  
 Supervised By : Jagrut Upadhyay 04/24/2023

Quant Time: Apr 20 01:48:24 2023  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_F\Methods\8270-BF041723.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Tue Apr 18 04:51:21 2023  
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc Units	Dev(Min)	
Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.763	152	239409	20.000 ng	0.00	
21) Naphthalene-d8	8.051	136	993494	20.000 ng	0.00	
39) Acenaphthene-d10	9.804	164	547314	20.000 ng	0.00	
64) Phenanthrene-d10	11.286	188	1016951	20.000 ng	0.00	
76) Chrysene-d12	13.921	240	642320	20.000 ng	0.00	
86) Perylene-d12	15.345	264	586284	20.000 ng	0.00	
System Monitoring Compounds						
5) 2-Fluorophenol	5.369	112	1724285	112.620 ng	0.00	
7) Phenol-d6	6.398	99	2262437	115.802 ng	0.00	
23) Nitrobenzene-d5	7.328	82	1396458	73.267 ng	0.00	
42) 2,4,6-Tribromophenol	10.592	330	681326	124.112 ng	0.00	
45) 2-Fluorobiphenyl	9.128	172	2439571	68.475 ng	0.00	
79) Terphenyl-d14	12.874	244	3107306	82.534 ng	0.00	
Target Compounds						
2) 1,4-Dioxane	2.493	88	342167	45.800 ng	99	Qvalue
3) Pyridine	3.199	79	859349	42.058 ng	98	
4) n-Nitrosodimethylamine	3.152	42	483407	49.792 ng	97	
6) Aniline	6.422	93	899615	34.722 ng	98	
8) 2-Chlorophenol	6.545	128	835394	53.783 ng	99	
9) Benzaldehyde	6.310	77	535106	46.291 ng	98	
10) Phenol	6.410	94	1110900	50.571 ng	99	
11) bis(2-Chloroethyl)ether	6.498	93	819427	49.392 ng	99	
12) 1,3-Dichlorobenzene	6.704	146	865037	50.668 ng	99	
13) 1,4-Dichlorobenzene	6.781	146	871175	50.925 ng	98	
14) 1,2-Dichlorobenzene	6.934	146	834738	52.261 ng	99	
15) Benzyl Alcohol	6.910	79	737914	55.112 ng	97	
16) 2,2'-oxybis(1-Chloropr...	7.045	45	1500972	52.287 ng	99	
17) 2-Methylphenol	7.022	107	701199	49.288 ng	99	
18) Hexachloroethane	7.281	117	303485	51.118 ng	94	
19) n-Nitroso-di-n-propyla...	7.187	70	615710	50.489 ng	100	
20) 3+4-Methylphenols	7.175	107	851198	50.621 ng	88	
22) Acetophenone	7.175	105	1106194	47.463 ng	95	
24) Nitrobenzene	7.351	77	921690	47.179 ng	95	
25) Isophorone	7.587	82	1739541	48.116 ng	99	
26) 2-Nitrophenol	7.663	139	447798	60.761 ng	95	
27) 2,4-Dimethylphenol	7.704	122	779365	51.842 ng	99	
28) bis(2-Chloroethoxy)met...	7.804	93	1026399	47.484 ng	100	
29) 2,4-Dichlorophenol	7.904	162	706866	50.516 ng	96	
30) 1,2,4-Trichlorobenzene	7.992	180	720441	47.683 ng	100	
31) Naphthalene	8.075	128	2336694	46.543 ng	99	
32) Benzoic acid	7.828	122	496238	65.824 ng	98	
33) 4-Chloroaniline	8.116	127	211152	10.269 ng	98	
34) Hexachlorobutadiene	8.192	225	396586	45.882 ng	100	
35) Caprolactam	8.492	113	260541m	62.523 ng		
36) 4-Chloro-3-methylphenol	8.598	107	782840	53.301 ng	98	
37) 2-Methylnaphthalene	8.763	142	1572829	45.515 ng	100	
38) 1-Methylnaphthalene	8.863	142	1487333	46.193 ng	100	
40) 1,2,4,5-Tetrachloroben...	8.933	216	671986	43.408 ng	98	
41) Hexachlorocyclopentadiene	8.922	237	871134	106.763 ng	98	
43) 2,4,6-Trichlorophenol	9.039	196	494381	50.379 ng	99	

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.075	196	532769	46.429	ng	99
46) 1,1'-Biphenyl	9.228	154	1960189	45.178	ng	99
47) 2-Chloronaphthalene	9.251	162	1487526	46.330	ng	98
48) 2-Nitroaniline	9.345	65	509334	54.835	ng	96
49) Acenaphthylene	9.669	152	2322634	45.355	ng	99
50) Dimethylphthalate	9.533	163	1817760	48.039	ng	100
51) 2,6-Dinitrotoluene	9.586	165	426798	51.445	ng	92
52) Acenaphthene	9.839	154	1698043	51.456	ng	99
53) 3-Nitroaniline	9.751	138	285854	30.279	ng	97
54) 2,4-Dinitrophenol	9.863	184	454631	140.675	ng #	86
55) Dibenzofuran	10.010	168	2159325	46.120	ng	100
56) 4-Nitrophenol	9.916	139	764519	109.017	ng	93
57) 2,4-Dinitrotoluene	9.992	165	573009	55.963	ng	96
58) Fluorene	10.351	166	1600633	46.670	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.128	232	462244	51.904	ng	100
60) Diethylphthalate	10.233	149	1768144	50.006	ng	99
61) 4-Chlorophenyl-phenyle...	10.345	204	796124	45.637	ng	93
62) 4-Nitroaniline	10.375	138	477358	52.607	ng	98
63) Azobenzene	10.504	77	1958622	49.589	ng	98
65) 4,6-Dinitro-2-methylph...	10.398	198	303682	62.354	ng	96
66) n-Nitrosodiphenylamine	10.463	169	1503511	45.395	ng	99
67) 4-Bromophenyl-phenylether	10.833	248	540789	46.569	ng	96
68) Hexachlorobenzene	10.904	284	589616	48.940	ng	97
69) Atrazine	10.992	200	536662	57.971	ng	99
70) Pentachlorophenol	11.092	266	727009	93.514	ng	99
71) Phenanthrene	11.310	178	2501933	45.738	ng	99
72) Anthracene	11.363	178	2538810	45.339	ng	99
73) Carbazole	11.516	167	2284627	47.153	ng	99
74) Di-n-butylphthalate	11.857	149	2821915	54.330	ng	99
75) Fluoranthene	12.498	202	2541389	46.719	ng	99
77) Benzidine	12.621	184	948810	94.826	ng	95
78) Pyrene	12.727	202	2609317	49.681	ng	99
80) Butylbenzylphthalate	13.351	149	927780	70.505	ng	91
81) Benzo(a)anthracene	13.910	228	2112462	48.359	ng	99
82) 3,3'-Dichlorobenzidine	13.874	252	635360	60.129	ng	99
83) Chrysene	13.951	228	2044460	47.088	ng	100
84) Bis(2-ethylhexyl)phtha...	13.916	149	1324531	79.725	ng	99
85) Di-n-octyl phthalate	14.527	149	2475873	65.829	ng	95
87) Indeno(1,2,3-cd)pyrene	16.745	276	1964922	58.653	ng	100
88) Benzo(b)fluoranthene	14.939	252	1963178	53.756	ng	99
89) Benzo(k)fluoranthene	14.968	252	1691089	46.485	ng	100
90) Benzo(a)pyrene	15.292	252	1549090	46.430	ng	99
91) Dibenzo(a,h)anthracene	16.768	278	1648387	57.748	ng	98
92) Benzo(g,h,i)perylene	17.168	276	1605849	58.757	ng	98

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

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