

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF061923\
 Data File : BF133833.D
 Acq On : 19 Jun 2023 14:40
 Operator : CG\JU
 Sample : PB153490BS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB153490BS

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 06/20/2023
 Supervised By :mohammad ahmed 06/20/2023

Quant Time: Jun 20 01:32:22 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF060923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 19 16:00:59 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.857	152	109009	20.000	ng	0.00
21) Naphthalene-d8	8.140	136	433162	20.000	ng	0.00
39) Acenaphthene-d10	9.898	164	224426	20.000	ng	0.00
64) Phenanthrene-d10	11.392	188	445244	20.000	ng	0.00
76) Chrysene-d12	14.051	240	202203	20.000	ng	0.00
86) Perylene-d12	15.557	264	206462	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.504	112	743891	118.801	ng	0.01
7) Phenol-d6	6.498	99	916009	112.386	ng	0.00
23) Nitrobenzene-d5	7.428	82	631971	79.465	ng	0.00
42) 2,4,6-Tribromophenol	10.698	330	316772	111.507	ng	0.00
45) 2-Fluorobiphenyl	9.222	172	1176879	74.544	ng	0.00
79) Terphenyl-d14	12.986	244	1282740	98.161	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.734	88	110506	40.142	ng	97
3) Pyridine	3.510	79	232443	28.969	ng	97
4) n-Nitrosodimethylamine	3.469	42	185217	41.760	ng	96
6) Aniline	6.522	93	292640	28.506	ng	# 89
8) 2-Chlorophenol	6.646	128	306277	46.414	ng	97
9) Benzaldehyde	6.410	77	84033	15.540	ng	96
10) Phenol	6.516	94	362869	42.636	ng	88
11) bis(2-Chloroethyl)ether	6.598	93	272453	43.121	ng	97
12) 1,3-Dichlorobenzene	6.798	146	317439	45.024	ng	98
13) 1,4-Dichlorobenzene	6.875	146	320346	44.955	ng	96
14) 1,2-Dichlorobenzene	7.028	146	307323	47.728	ng	99
15) Benzyl Alcohol	7.004	79	276583	43.448	ng	99
16) 2,2'-oxybis(1-Chloropr...	7.134	45	494780	46.196	ng	99
17) 2-Methylphenol	7.116	107	253100	43.273	ng	99
18) Hexachloroethane	7.369	117	120911	49.563	ng	99
19) n-Nitroso-di-n-propyla...	7.275	70	220736	42.566	ng	96
20) 3+4-Methylphenols	7.269	107	302610	39.772	ng	# 84
22) Acetophenone	7.269	105	427079	42.988	ng	92
24) Nitrobenzene	7.445	77	333220	41.611	ng	99
25) Isophorone	7.681	82	605213	41.452	ng	99
26) 2-Nitrophenol	7.757	139	164346	45.729	ng	97
27) 2,4-Dimethylphenol	7.793	122	271428	43.818	ng	99
28) bis(2-Chloroethoxy)met...	7.893	93	350430	41.528	ng	99
29) 2,4-Dichlorophenol	7.998	162	258214	42.532	ng	99
30) 1,2,4-Trichlorobenzene	8.081	180	266093	42.010	ng	99
31) Naphthalene	8.163	128	849915	40.273	ng	99
32) Benzoic acid	7.922	122	207267	53.054	ng	94
33) 4-Chloroaniline	8.210	127	99242	10.998	ng	97
34) Hexachlorobutadiene	8.275	225	169071	40.769	ng	98
35) Caprolactam	8.592	113	87387m	42.777	ng	
36) 4-Chloro-3-methylphenol	8.698	107	298000	42.585	ng	99
37) 2-Methylnaphthalene	8.857	142	579311	39.508	ng	99
38) 1-Methylnaphthalene	8.951	142	545146	40.402	ng	99
40) 1,2,4,5-Tetrachloroben...	9.022	216	274000	40.160	ng	# 98
41) Hexachlorocyclopentadiene	9.004	237	357867	117.703	ng	98
43) 2,4,6-Trichlorophenol	9.134	196	190913	41.874	ng	97

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
44) 2,4,5-Trichlorophenol	9.175	196	204025	38.465	ng	97
46) 1,1'-Biphenyl	9.322	154	746363	41.235	ng	99
47) 2-Chloronaphthalene	9.345	162	544425	41.932	ng	99
48) 2-Nitroaniline	9.445	65	195881	41.252	ng	99
49) Acenaphthylene	9.763	152	912371	40.374	ng	100
50) Dimethylphthalate	9.628	163	675923	40.350	ng	100
51) 2,6-Dinitrotoluene	9.687	165	146578	42.392	ng	# 76
52) Acenaphthene	9.934	154	535222	40.463	ng	100
53) 3-Nitroaniline	9.857	138	94589	22.480	ng	93
54) 2,4-Dinitrophenol	9.969	184	182735	106.139	ng	96
55) Dibenzofuran	10.110	168	820568	40.655	ng	98
56) 4-Nitrophenol	10.028	139	242762	85.486	ng	86
57) 2,4-Dinitrotoluene	10.092	165	193778	44.066	ng	92
58) Fluorene	10.451	166	648589	42.020	ng	100
59) 2,3,4,6-Tetrachlorophenol	10.228	232	180178	45.372	ng	97
60) Diethylphthalate	10.328	149	713925	41.875	ng	99
61) 4-Chlorophenyl-phenyle...	10.445	204	307873	40.404	ng	98
62) 4-Nitroaniline	10.481	138	160844	38.987	ng	96
63) Azobenzene	10.604	77	666252	41.280	ng	96
65) 4,6-Dinitro-2-methylph...	10.504	198	116076	54.557	ng	83
66) n-Nitrosodiphenylamine	10.563	169	576583	38.552	ng	99
67) 4-Bromophenyl-phenylether	10.933	248	196755	39.592	ng	95
68) Hexachlorobenzene	10.998	284	214487	41.044	ng	96
69) Atrazine	11.092	200	188142	43.357	ng	97
70) Pentachlorophenol	11.198	266	237340	75.845	ng	97
71) Phenanthrene	11.422	178	912609	40.400	ng	99
72) Anthracene	11.469	178	946789	40.197	ng	99
73) Carbazole	11.628	167	862573	39.320	ng	99
74) Di-n-butylphthalate	11.957	149	1108656	39.421	ng	99
75) Fluoranthene	12.610	202	942809	39.303	ng	98
77) Benzidine	12.733	184	220010	32.425	ng	99
78) Pyrene	12.845	202	937980	49.797	ng	100
80) Butylbenzylphthalate	13.469	149	350772	44.112	ng	96
81) Benzo(a)anthracene	14.039	228	583628	41.728	ng	99
82) 3,3'-Dichlorobenzidine	14.004	252	138416	29.767	ng	97
83) Chrysene	14.074	228	549748	41.557	ng	99
84) Bis(2-ethylhexyl)phtha...	14.027	149	396612	40.507	ng	99
85) Di-n-octyl phthalate	14.651	149	572659	41.158	ng	98
87) Indeno(1,2,3-cd)pyrene	17.110	276	706836	49.763	ng	100
88) Benzo(b)fluoranthene	15.110	252	504158	40.199	ng	99
89) Benzo(k)fluoranthene	15.139	252	512369	41.939	ng	99
90) Benzo(a)pyrene	15.492	252	480564	41.287	ng	98
91) Dibenzo(a,h)anthracene	17.133	278	576125	48.993	ng	99
92) Benzo(g,h,i)perylene	17.586	276	577580	49.539	ng	99

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

