

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF101822\
 Data File : BF130705.D
 Acq On : 18 Oct 2022 18:52
 Operator : CG\JU
 Sample : N5172-01MS
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BU-02-101722MS

Quant Time: Oct 19 02:40:34 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

Manual Integrations APPROVED

Reviewed By : Christian Giraldo 10/19/2022
 Supervised By : Jagrut Upadhyay 10/19/2022

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.575	152	58230	20.000	ng	0.00
21) Naphthalene-d8	7.863	136	194208	20.000	ng	0.00
39) Acenaphthene-d10	9.610	164	91388	20.000	ng	0.00
64) Phenanthrene-d10	11.092	188	166763	20.000	ng	0.00
76) Chrysene-d12	13.745	240	140729	20.000	ng	0.02
86) Perylene-d12	15.115	264	102988	20.000	ng	# 0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.187	112	385148	119.225	ng	0.02
7) Phenol-d6	6.234	99	429501	105.748	ng	0.00
23) Nitrobenzene-d5	7.151	82	262502	73.788	ng	0.00
42) 2,4,6-Tribromophenol	10.404	330	113366	126.776	ng	0.00
45) 2-Fluorobiphenyl	8.933	172	442312	73.313	ng	0.00
79) Terphenyl-d14	12.686	244	588418	69.397	ng	0.01
Target Compounds						
						Qvalue
2) 1,4-Dioxane	2.169	88	64195	32.847	ng	97
3) Pyridine	2.834	79	158376	45.303	ng	96
4) n-Nitrosodimethylamine	2.793	42	80839	49.078	ng	93
6) Aniline	6.239	93	162242	33.365	ng	94
8) 2-Chlorophenol	6.369	128	167324	44.044	ng	94
9) Benzaldehyde	6.128	77	79771	31.261	ng	96
10) Phenol	6.245	94	196977	41.915	ng	99
11) bis(2-Chloroethyl)ether	6.322	93	161078	47.186	ng	97
12) 1,3-Dichlorobenzene	6.516	146	188649	45.353	ng	98
13) 1,4-Dichlorobenzene	6.592	146	192032	45.507	ng	99
14) 1,2-Dichlorobenzene	6.745	146	178396	45.393	ng	97
15) Benzyl Alcohol	6.734	79	131378	45.054	ng	98
16) 2,2'-oxybis(1-Chloropr...	6.863	45	184132	41.926	ng	# 49
17) 2-Methylphenol	6.857	107	126186	41.757	ng	92
18) Hexachloroethane	7.081	117	66296	41.528	ng	96
19) n-Nitroso-di-n-propyla...	7.004	70	110606	42.209	ng	95
20) 3+4-Methylphenols	7.016	107	162664	40.170	ng	# 79
22) Acetophenone	6.998	105	235174	50.480	ng	# 93
24) Nitrobenzene	7.169	77	157918	44.140	ng	97
25) Isophorone	7.410	82	277062	46.793	ng	99
26) 2-Nitrophenol	7.486	139	80814	46.061	ng	97
27) 2,4-Dimethylphenol	7.539	122	130955	53.553	ng	97
28) bis(2-Chloroethoxy)met...	7.622	93	167212	48.773	ng	99
29) 2,4-Dichlorophenol	7.739	162	116356	42.529	ng	97
30) 1,2,4-Trichlorobenzene	7.804	180	129575	43.931	ng	97
31) Naphthalene	7.886	128	1744844	173.563	ng	99
32) Benzoic acid	7.681	122	67165	43.416	ng	95
33) 4-Chloroaniline	7.945	127	103925	26.686	ng	98
34) Hexachlorobutadiene	7.998	225	85235	46.560	ng	98
35) Caprolactam	8.316	113	36446	44.276	ng	98
36) 4-Chloro-3-methylphenol	8.439	107	113751	38.330	ng	94
37) 2-Methylnaphthalene	8.575	142	833518	131.752	ng	99
38) 1-Methylnaphthalene	8.675	142	683920	111.360	ng	100
40) 1,2,4,5-Tetrachloroben...	8.739	216	117010	47.192	ng	99
41) Hexachlorocyclopentadiene	8.722	237	5171	14.738	ng	94
43) 2,4,6-Trichlorophenol	8.863	196	72730	43.687	ng	99

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44) 2,4,5-Trichlorophenol	8.910	196	78819	43.051	ng	98
46) 1,1'-Biphenyl	9.033	154	432691	65.331	ng	99
47) 2-Chloronaphthalene	9.057	162	235621	43.529	ng	98
48) 2-Nitroaniline	9.169	65	70603	43.329	ng	92
49) Acenaphthylene	9.475	152	584950	72.123	ng	99
50) Dimethylphthalate	9.339	163	273382	43.215	ng	98
51) 2,6-Dinitrotoluene	9.410	165	58360	42.056	ng	95
52) Acenaphthene	9.645	154	402321	81.875	ng	99
53) 3-Nitroaniline	9.580	138	53392	35.100	ng	96
54) 2,4-Dinitrophenol	9.716	184	10870	19.390	ng	# 19
55) Dibenzofuran	9.816	168	820130	109.559	ng	98
56) 4-Nitrophenol	9.775	139	73800	87.992	ng	98
57) 2,4-Dinitrotoluene	9.822	165	82603	46.316	ng	# 89
58) Fluorene	10.163	166	1129977	183.609	ng	99
59) 2,3,4,6-Tetrachlorophenol	9.945	232	61316	42.008	ng	# 75
60) Diethylphthalate	10.039	149	264560	41.915	ng	99
61) 4-Chlorophenyl-phenyle...	10.151	204	119930	43.104	ng	96
62) 4-Nitroaniline	10.192	138	56750	39.076	ng	96
63) Azobenzene	10.310	77	236909	39.850	ng	# 72
65) 4,6-Dinitro-2-methylph...	10.239	198	10836m	10.398	ng	
66) n-Nitrosodiphenylamine	10.275	169	233659m	42.122	ng	
67) 4-Bromophenyl-phenylether	10.639	248	80258	42.534	ng	98
68) Hexachlorobenzene	10.710	284	94309	44.214	ng	96
69) Atrazine	10.816	200	77227	51.126	ng	97
70) Pentachlorophenol	10.922	266	79930	96.282	ng	97
71) Phenanthrene	11.139	178	4877080	527.865	ng	97
72) Anthracene	11.180	178	1266607	140.368	ng	99
73) Carbazole	11.339	167	811518	100.788	ng	100
74) Di-n-butylphthalate	11.663	149	465218	47.698	ng	99
75) Fluoranthene	12.333	202	4127814	464.851	ng	96
77) Benzidine	12.457	184	131413	36.944	ng	97
78) Pyrene	12.563	202	3826708	315.195	ng	97
80) Butylbenzylphthalate	13.163	149	231820	49.657	ng	98
81) Benzo(a)anthracene	13.733	228	1950854	205.732	ng	96
82) 3,3'-Dichlorobenzidine	13.698	252	112917	41.208	ng	100
83) Chrysene	13.774	228	1565067	174.077	ng	97
84) Bis(2-ethylhexyl)phtha...	13.721	149	330122	58.328	ng	99
85) Di-n-octyl phthalate	14.327	149	483283	55.756	ng	97
87) Indeno(1,2,3-cd)pyrene	16.427	276	705846	91.570	ng	96
88) Benzo(b)fluoranthene	14.745	252	1422156	218.075	ng	96
89) Benzo(k)fluoranthene	14.768	252	656465m	98.861	ng	
90) Benzo(a)pyrene	15.074	252	1094934	201.732	ng	97
91) Dibenzo(a,h)anthracene	16.433	278	352386	55.771	ng	97
92) Benzo(g,h,i)perylene	16.821	276	660092	112.167	ng	97

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

