

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011720\
 Data File : BF118562.D
 Acq On : 17 Jan 2020 16:58
 Operator : JU
 Sample : PB126158BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB126158BS

Manual Integrations
 APPROVED

mohammad
 1/21/2020 7:58:31 AM

Quant Time: Jan 18 01:20:54 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF011720.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 17 15:49:53 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	431857	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	1640978	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	950582	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	1725785	20.00	ng	0.00
76) Chrysene-d12	14.05	240	1282541	20.00	ng	0.00
87) Perylene-d12	15.53	264	1088392	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	2868791	122.74	ng	0.02
7) Phenol-d6	6.52	99	3770618	123.12	ng	0.00
23) Nitrobenzene-d5	7.45	82	2216565	83.43	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	1342134	127.72	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	4242928	80.58	ng	0.00
79) Terphenyl-d14	13.00	244	4688092	72.19	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	2.78	88	523326	46.600	ng	99
3) Pyridine	3.53	79	1134112	36.566	ng	100
4) n-Nitrosodimethylamine	3.47	42	586077	41.198	ng	99
6) Aniline	6.55	93	1384000	34.527	ng	99
8) 2-Chlorophenol	6.68	128	1216631	45.770	ng	95
9) Benzaldehyde	6.43	77	663672	42.334	ng	99
10) Phenol	6.53	94	1535774	45.707	ng	99
11) bis(2-Chloroethyl)ether	6.62	93	1148100	42.662	ng	96
12) 1,3-Dichlorobenzene	6.83	146	1308202	41.386	ng	99
13) 1,4-Dichlorobenzene	6.91	146	1328197	41.825	ng	98
14) 1,2-Dichlorobenzene	7.06	146	1263434	42.524	ng	97
15) Benzyl Alcohol	7.03	79	1123078	44.685	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.16	45	1597977	40.477	ng	99
17) 2-Methylphenol	7.14	107	1019855	45.200	ng	99
18) Hexachloroethane	7.40	117	480123	41.489	ng	97
19) n-Nitroso-di-n-propylamine	7.30	70	943592	42.965	ng	99
20) 3+4-Methylphenols	7.29	107	1263492	46.828	ng	# 89
22) Acetophenone	7.30	105	1661071	41.713	ng	# 95
24) Nitrobenzene	7.47	77	1312367	46.447	ng	100
25) Isophorone	7.71	82	2472937	42.885	ng	98
26) 2-Nitrophenol	7.79	139	587589	48.099	ng	95
27) 2,4-Dimethylphenol	7.82	122	1178946	51.407	ng	99
28) bis(2-Chloroethoxy)methane	7.92	93	1518039	40.973	ng	100
29) 2,4-Dichlorophenol	8.03	162	1107761	43.883	ng	99
30) 1,2,4-Trichlorobenzene	8.12	180	1198210	40.730	ng	99
31) Naphthalene	8.20	128	3396667	43.891	ng	99
32) Benzoic acid	7.94	122	699076	45.070	ng	99
33) 4-Chloroaniline	8.23	127	824448	22.928	ng	99
34) Hexachlorobutadiene	8.32	225	710921	38.743	ng	98
35) Caprolactam	8.61	113	345673m	44.018	ng	
36) 4-Chloro-3-methylphenol	8.72	107	1118507	44.697	ng	96
37) 2-Methylnaphthalene	8.89	142	2449414	44.676	ng	100
38) 1-Methylnaphthalene	8.99	142	2327627	44.876	ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.06	216	1218127	39.730	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.05	237	1562993	97.568	nq	98
43) 2,4,6-Trichlorophenol	9.16	196	857600	43.901	nq	98
44) 2,4,5-Trichlorophenol	9.20	196	885662	44.843	nq	98
46) 1,1'-Biphenyl	9.35	154	2910426	42.849	nq	99
47) 2-Chloronaphthalene	9.37	162	2356964	41.170	nq	98
48) 2-Nitroaniline	9.46	65	696683	46.789	nq	98
49) Acenaphthylene	9.79	152	3555753	43.989	nq	100
50) Dimethylphthalate	9.65	163	2712459	41.560	nq	100
51) 2,6-Dinitrotoluene	9.71	165	628368	50.460	nq	94
52) Acenaphthene	9.96	154	2407110	45.457	nq	100
53) 3-Nitroaniline	9.87	138	400053	28.359	nq	# 91
54) 2,4-Dinitrophenol	9.98	184	480924	89.136	nq	# 1
55) Dibenzofuran	10.14	168	3246046	43.317	nq	98
56) 4-Nitrophenol	10.03	139	1065953	102.291	nq	94
57) 2,4-Dinitrotoluene	10.11	165	815094	49.485	nq	99
58) Fluorene	10.48	166	2504791	43.164	nq	100
59) 2,3,4,6-Tetrachlorophenol	10.25	232	787611	46.615	nq	99
60) Diethylphthalate	10.35	149	2685488	42.147	nq	99
61) 4-Chlorophenyl-phenylether	10.47	204	1373846	43.275	nq	99
62) 4-Nitroaniline	10.49	138	612100	43.687	nq	99
63) Azobenzene	10.63	77	2666509	42.938	nq	100
65) 4,6-Dinitro-2-methylphenol	10.52	198	351451	49.625	nq	98
66) n-Nitrosodiphenylamine	10.59	169	2350959	42.396	nq	99
67) 4-Bromophenyl-phenylether	10.96	248	943399	40.837	nq	99
68) Hexachlorobenzene	11.03	284	1015501	39.668	nq	99
69) Atrazine	11.12	200	853208	53.318	nq	98
70) Pentachlorophenol	11.22	266	1171418	93.107	nq	99
71) Phenanthrene	11.44	178	3606155	41.571	nq	100
72) Anthracene	11.49	178	3705794	42.818	nq	99
73) Carbazole	11.64	167	3282661	42.055	nq	99
74) Di-n-butylphthalate	11.97	149	3815005	39.330	nq	99
75) Fluoranthene	12.63	202	3796848	40.701	nq	99
77) Benzidine	12.74	184	1966147	63.096	nq	98
78) Pyrene	12.86	202	3818530	39.891	nq	99
80) Butylbenzylphthalate	13.47	149	1669991	40.699	nq	99
81) Benzo(a)anthracene	14.04	228	3315493	40.955	nq	99
82) 3,3'-Dichlorobenzidine	14.00	252	856645	30.161	nq	100
83) Chrysene	14.08	228	3222739	41.387	nq	99
84) Bis(2-ethylhexyl)phthalate	14.04	149	2148420	42.184	nq	# 98
85) Di-n-octyl phthalate	14.65	149	3386957	43.427	nq	99
86) Indeno(1,2,3-cd)pyrene	17.01	276	2762777	36.055	nq	100
88) Benzo(b)fluoranthene	15.10	252	3147051	44.395	nq	100
89) Benzo(k)fluoranthene	15.13	252	2663873	41.580	nq	99
90) Benzo(a)pyrene	15.47	252	2533640	41.076	nq	99
91) Dibenzo(a,h)anthracene	17.03	278	2318729	40.249	nq	100
92) Benzo(g,h,i)perylene	17.46	276	2244559	40.336	nq	99

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

