

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080519\
 Data File : BF115982.D
 Acq On : 5 Aug 2019 16:24
 Operator : HP/JU
 Sample : PB121965BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB121965BS

Manual Integrations
 APPROVED

mohammad
 8/8/2019 4:45:43 PM

Quant Time: Aug 06 02:03:45 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 31 15:31:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	133944	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	531846	20.00	ng	-0.01
39) Acenaphthene-d10	9.88	164	275113	20.00	ng	-0.02
64) Phenanthrene-d10	11.36	188	452314	20.00	ng	-0.02
76) Chrysene-d12	14.00	240	307898	20.00	ng	-0.02
87) Perylene-d12	15.47	264	285202	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.48	112	823570	102.93	ng	0.01
7) Phenol-d6	6.48	99	1041260	103.15	ng	0.00
23) Nitrobenzene-d5	7.40	82	679211	73.72	ng	-0.01
42) 2,4,6-Tribromophenol	10.67	330	276921	103.82	ng	-0.02
45) 2-Fluorobiphenyl	9.20	172	1210153	82.39	ng	-0.02
79) Terphenyl-d14	12.95	244	1219333	83.45	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.71	88	153743	38.036	ng	100
3) Pyridine	3.46	79	332846	29.478	ng	99
4) n-Nitrosodimethylamine	3.40	42	151161	33.923	ng	97
6) Aniline	6.51	93	376336	26.031	ng	99
8) 2-Chlorophenol	6.63	128	338173	38.222	ng	99
9) Benzaldehyde	6.40	77	124308	17.847	ng	99
10) Phenol	6.49	94	429576	36.136	ng	99
11) bis(2-Chloroethyl)ether	6.58	93	327159	37.433	ng	97
12) 1,3-Dichlorobenzene	6.79	146	365407	35.736	ng	98
13) 1,4-Dichlorobenzene	6.86	146	373738	36.505	ng	99
14) 1,2-Dichlorobenzene	7.02	146	348579	36.976	ng	99
15) Benzyl Alcohol	6.99	79	280780	36.651	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.12	45	439396	36.858	ng	95
17) 2-Methylphenol	7.10	107	280588	37.453	ng	99
18) Hexachloroethane	7.36	117	134675	35.728	ng	97
19) n-Nitroso-di-n-propylamine	7.26	70	233676	37.355	ng	98
20) 3+4-Methylphenols	7.25	107	343101	38.701	ng	# 80
22) Acetophenone	7.25	105	466354	39.982	ng	97
24) Nitrobenzene	7.42	77	376297	37.824	ng	99
25) Isophorone	7.66	82	668239	37.929	ng	100
26) 2-Nitrophenol	7.74	139	183110	39.046	ng	98
27) 2,4-Dimethylphenol	7.78	122	299450	42.442	ng	99
28) bis(2-Chloroethoxy)methane	7.87	93	414814	37.987	ng	99
29) 2,4-Dichlorophenol	7.99	162	288170	38.682	ng	99
30) 1,2,4-Trichlorobenzene	8.07	180	320309	37.719	ng	100
31) Naphthalene	8.15	128	986929	39.434	ng	100
32) Benzoic acid	7.91	122	170289	34.234	ng	98
33) 4-Chloroaniline	8.19	127	270144	24.158	ng	98
34) Hexachlorobutadiene	8.26	225	178936	36.583	ng	99
35) Caprolactam	8.56	113	83566m	34.683	ng	
36) 4-Chloro-3-methylphenol	8.68	107	295914	38.182	ng	96
37) 2-Methylnaphthalene	8.84	142	651959	39.554	ng	99
38) 1-Methylnaphthalene	8.94	142	615248	39.456	ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.00	216	297289	40.421	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.99	237	170380	53.704	nq	100
43) 2,4,6-Trichlorophenol	9.12	196	184427	36.430	nq	99
44) 2,4,5-Trichlorophenol	9.16	196	197209	37.685	nq	96
46) 1,1'-Biphenyl	9.30	154	798318	41.102	nq	99
47) 2-Chloronaphthalene	9.33	162	614460	38.010	nq	99
48) 2-Nitroaniline	9.42	65	187827	35.152	nq	90
49) Acenaphthylene	9.74	152	934598	39.628	nq	99
50) Dimethylphthalate	9.60	163	707777	37.784	nq	100
51) 2,6-Dinitrotoluene	9.66	165	155886	38.294	nq	93
52) Acenaphthene	9.91	154	560842	38.763	nq	99
53) 3-Nitroaniline	9.83	138	117538	23.367	nq	94
54) 2,4-Dinitrophenol	9.95	184	121674	60.840	nq	90
55) Dibenzofuran	10.09	168	843507	40.089	nq	99
56) 4-Nitrophenol	10.00	139	212168	65.845	nq	99
57) 2,4-Dinitrotoluene	10.07	165	203884	37.617	nq	90
58) Fluorene	10.43	166	640073	39.539	nq	100
59) 2,3,4,6-Tetrachlorophenol	10.21	232	167661	38.082	nq	98
60) Diethylphthalate	10.30	149	676490	38.681	nq	100
61) 4-Chlorophenyl-phenylether	10.42	204	326571	39.832	nq	99
62) 4-Nitroaniline	10.45	138	159022	30.592	nq	99
63) Azobenzene	10.57	77	704684	39.175	nq	99
65) 4,6-Dinitro-2-methylphenol	10.49	198	92846	38.735	nq	93
66) n-Nitrosodiphenylamine	10.53	169	560328	41.158	nq	99
67) 4-Bromophenyl-phenylether	10.90	248	194938	40.260	nq	99
68) Hexachlorobenzene	10.98	284	216234	38.620	nq	98
69) Atrazine	11.06	200	168326	37.583	nq	99
70) Pentachlorophenol	11.17	266	189234	69.615	nq	97
71) Phenanthrene	11.39	178	904425	39.113	nq	100
72) Anthracene	11.44	178	938321	40.096	nq	99
73) Carbazole	11.59	167	832649	39.218	nq	99
74) Di-n-butylphthalate	11.92	149	1039838	41.985	nq	100
75) Fluoranthene	12.58	202	926790	38.285	nq	100
77) Benzidine	12.70	184	299557	28.798	nq	98
78) Pyrene	12.81	202	933545	40.222	nq	99
80) Butylbenzylphthalate	13.42	149	416042	42.376	nq	96
81) Benzo(a)anthracene	13.99	228	759856	38.611	nq	100
82) 3,3'-Dichlorobenzidine	13.95	252	243782	35.656	nq	98
83) Chrysene	14.03	228	738068	38.630	nq	99
84) Bis(2-ethylhexyl)phthalate	13.97	149	597227	46.811	nq	99
85) Di-n-octyl phthalate	14.59	149	933813	46.081	nq	100
86) Indeno(1,2,3-cd)pyrene	16.94	276	737711	45.972	nq	99
88) Benzo(b)fluoranthene	15.04	252	692965	42.021	nq	98
89) Benzo(k)fluoranthene	15.07	252	665155	41.411	nq	99
90) Benzo(a)pyrene	15.40	252	647078	43.895	nq	98
91) Dibenzo(a,h)anthracene	16.94	278	627584	49.183	nq	100
92) Benzo(g,h,i)perylene	17.38	276	618677	49.369	nq	98

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

