

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121019\
 Data File : BP001298.D
 Acq On : 10 Dec 2019 16:46
 Operator : JU
 Sample : PB125370BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB125370BS

Manual Integrations
 APPROVED

mohammad
 12/11/2019 12:31:01 PM

Quant Time: Dec 11 04:30:03 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP112719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 05 12:34:47 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.30	152	71525	20.00	ng	0.00
21) Naphthalene-d8	10.34	136	261833	20.00	ng	-0.01
39) Acenaphthene-d10	13.21	164	152338	20.00	ng	0.00
64) Phenanthrene-d10	15.61	188	335054	20.00	ng	0.00
76) Chrysene-d12	19.26	240	320315	20.00	ng	0.00
87) Perylene-d12	21.02	264	284076	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	6.21	112	495780	115.00	ng	0.00
7) Phenol-d6	7.76	99	666971	116.13	ng	0.00
23) Nitrobenzene-d5	9.19	82	465419	77.12	ng	-0.01
42) 2,4,6-Tribromophenol	14.50	330	304780	208.73	ng	0.00
45) 2-Fluorobiphenyl	12.12	172	942158	90.52	ng	-0.01
79) Terphenyl-d14	17.89	244	1291379	74.59	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.56	88	55506	27.564	ng	98
3) Pyridine	3.41	79	151799	27.166	ng	93
4) n-Nitrosodimethylamine	3.31	42	97137	40.173	ng	84
6) Aniline	7.78	93	193423	25.400	ng	# 1
8) 2-Chlorophenol	7.96	128	173242	38.840	ng	98
9) Benzaldehyde	7.59	77	63070	14.425	ng	99
10) Phenol	7.78	94	236213	36.157	ng	97
11) bis(2-Chloroethyl)ether	7.91	93	167213	32.869	ng	97
12) 1,3-Dichlorobenzene	8.21	146	192605	36.431	ng	99
13) 1,4-Dichlorobenzene	8.33	146	195895	36.783	ng	100
14) 1,2-Dichlorobenzene	8.57	146	188809	37.670	ng	96
15) Benzyl Alcohol	8.53	79	196169	41.608	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.77	45	259466	31.724	ng	99
17) 2-Methylphenol	8.73	107	150758	36.846	ng	99
18) Hexachloroethane	9.11	117	80093	36.354	ng	# 87
19) n-Nitroso-di-n-propylamine	8.97	70	150203	36.242	ng	98
20) 3+4-Methylphenols	8.97	107	201823	39.131	ng	98
22) Acetophenone	8.95	105	286292	37.202	ng	98
24) Nitrobenzene	9.22	77	228876	38.139	ng	99
25) Isophorone	9.61	82	404731	37.399	ng	99
26) 2-Nitrophenol	9.73	139	90252	41.059	ng	96
27) 2,4-Dimethylphenol	9.82	122	154404	44.346	ng	97
28) bis(2-Chloroethoxy)methane	9.98	93	228099	33.594	ng	98
29) 2,4-Dichlorophenol	10.12	162	170918	43.130	ng	96
30) 1,2,4-Trichlorobenzene	10.26	180	193293	41.761	ng	97
31) Naphthalene	10.37	128	511494	38.250	ng	99
32) Benzoic acid	10.01	122	95209	37.824	ng	91
33) 4-Chloroaniline	10.47	127	111189	19.161	ng	98
34) Hexachlorobutadiene	10.60	225	136123	46.751	ng	94
35) Caprolactam	11.04	113	52189m	38.212	ng	
36) 4-Chloro-3-methylphenol	11.29	107	200097	42.588	ng	94
37) 2-Methylnaphthalene	11.51	142	374756	41.070	ng	97
38) 1-Methylnaphthalene	11.67	142	356041	41.306	ng	100
40) 1,2,4,5-Tetrachlorobenzene	11.79	216	223606	46.576	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	11.78	237	261130	117.904	nq	100
43) 2,4,6-Trichlorophenol	11.97	196	141998	45.300	nq	99
44) 2,4,5-Trichlorophenol	12.03	196	155269	47.807	nq	96
46) 1,1'-Biphenyl	12.28	154	483046	41.029	nq	99
47) 2-Chloronaphthalene	12.30	162	380680	40.480	nq	98
48) 2-Nitroaniline	12.47	65	141804	38.474	nq	96
49) Acenaphthylene	12.97	152	600662	42.611	nq	99
50) Dimethylphthalate	12.80	163	489971	43.005	nq	99
51) 2,6-Dinitrotoluene	12.88	165	106729	43.760	nq	99
52) Acenaphthene	13.26	154	345136	40.703	nq	99
53) 3-Nitroaniline	13.14	138	68531	25.014	nq	94
54) 2,4-Dinitrophenol	13.32	184	87897	87.500	nq	94
55) Dibenzofuran	13.55	168	565948	44.416	nq	98
56) 4-Nitrophenol	13.45	139	155674	80.187	nq	91
57) 2,4-Dinitrotoluene	13.54	165	144560	47.287	nq	96
58) Fluorene	14.11	166	456576	44.718	nq	99
59) 2,3,4,6-Tetrachlorophenol	13.76	232	145878	54.640	nq	93
60) Diethylphthalate	13.97	149	506131	42.903	nq	99
61) 4-Chlorophenyl-phenylether	14.13	204	248211	47.740	nq	95
62) 4-Nitroaniline	12.47	138	119023	37.471	nq	97
63) Azobenzene	14.39	77	500706	39.265	nq	96
65) 4,6-Dinitro-2-methylphenol	14.21	198	62143	43.867	nq	94
66) n-Nitrosodiphenylamine	14.32	169	399031	39.196	nq	100
67) 4-Bromophenyl-phenylether	14.93	248	158783	43.651	nq	91
68) Hexachlorobenzene	15.01	284	184393	46.109	nq	# 90
69) Atrazine	15.22	200	138381	44.519	nq	98
70) Pentachlorophenol	15.33	266	319859	153.460	nq	94
71) Phenanthrene	15.65	178	700631	39.775	nq	99
72) Anthracene	15.72	178	717277	41.313	nq	98
73) Carbazole	15.97	167	620738	39.290	nq	99
74) Di-n-butylphthalate	16.52	149	821572	38.677	nq	99
75) Fluoranthene	17.36	202	806719	43.260	nq	97
77) Benzidine	17.54	184	237447	28.711	nq	99
78) Pyrene	17.66	202	814835	32.298	nq	99
80) Butylbenzylphthalate	18.55	149	347483	29.820	nq	93
81) Benzo(a)anthracene	19.24	228	766780	34.526	nq	100
82) 3,3'-Dichlorobenzidine	19.21	252	196628	26.800	nq	# 96
83) Chrysene	19.29	228	702531	35.326	nq	99
84) Bis(2-ethylhexyl)phthalate	19.29	149	492196	30.722	nq	# 98
85) Di-n-octyl phthalate	20.07	149	810866	28.492	nq	98
86) Indeno(1,2,3-cd)pyrene	22.67	276	773161	32.988	nq	# 94
88) Benzo(b)fluoranthene	20.54	252	721459	41.580	nq	98
89) Benzo(k)fluoranthene	20.57	252	693334	41.487	nq	99
90) Benzo(a)pyrene	20.96	252	672258	42.063	nq	# 97
91) Dibenzo(a,h)anthracene	22.70	278	653004	41.751	nq	# 96
92) Benzo(g,h,i)perylene	23.16	276	628249	40.679	nq	# 94

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

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