

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP062620\
 Data File : BP002551.D
 Acq On : 26 Jun 2020 15:00
 Operator : CG/JU
 Sample : PB129799BS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB129799BS

Manual Integrations
 APPROVED

mohammad
 6/29/2020 3:18:47 PM

Quant Time: Jun 26 17:43:13 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP060520.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 24 13:54:06 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	192824	20.00	ng	0.00
21) Naphthalene-d8	10.36	136	732735	20.00	ng	0.00
39) Acenaphthene-d10	14.23	164	437645	20.00	ng	0.00
64) Phenanthrene-d10	16.99	188	890983	20.00	ng	0.00
76) Chrysene-d12	21.10	240	696620	20.00	ng	0.00
86) Perylene-d12	23.29	264	646864	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.20	112	899409	86.31	ng	0.00
7) Phenol-d6	6.78	99	1134605	81.78	ng	0.00
23) Nitrobenzene-d5	8.73	82	1106511	90.18	ng	0.00
42) 2,4,6-Tribromophenol	15.73	330	374315	89.33	ng	0.00
45) 2-Fluorobiphenyl	12.85	172	2248127	86.37	ng	0.00
79) Terphenyl-d14	19.58	244	2745346	109.20	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.16	88	151736	31.611	ng	99
3) Pyridine	3.53	79	447687	34.315	ng	99
4) n-Nitrosodimethylamine	3.45	42	227831	42.376	ng	98
6) Aniline	6.92	93	602244	31.918	ng	99
8) 2-Chlorophenol	7.15	128	510665	43.584	ng	100
9) Benzaldehyde	6.73	77	224955	27.845	ng	99
10) Phenol	6.80	94	646971	42.996	ng	98
11) bis(2-Chloroethyl)ether	7.02	93	488940	38.919	ng	98
12) 1,3-Dichlorobenzene	7.48	146	544107	40.328	ng	99
13) 1,4-Dichlorobenzene	7.62	146	553535	40.778	ng	99
14) 1,2-Dichlorobenzene	7.93	146	530415	40.791	ng	98
15) Benzyl Alcohol	7.82	79	438844	40.810	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.12	45	680630	38.183	ng	99
17) 2-Methylphenol	8.03	107	430574	39.160	ng	98
18) Hexachloroethane	8.65	117	204897	41.433	ng	100
19) n-Nitroso-di-n-propylamine	8.39	70	370153	39.915	ng	100
20) 3+4-Methylphenols	8.36	107	576974	38.882	ng	97
22) Acetophenone	8.40	105	710029	43.131	ng	100
24) Nitrobenzene	8.77	77	584831	44.342	ng	99
25) Isophorone	9.29	82	1059439	42.396	ng	100
26) 2-Nitrophenol	9.48	139	280148	47.667	ng	98
27) 2,4-Dimethylphenol	9.55	122	443750	48.129	ng	98
28) bis(2-Chloroethoxy)methane	9.78	93	652103	43.811	ng	100
29) 2,4-Dichlorophenol	10.02	162	483500	46.585	ng	98
30) 1,2,4-Trichlorobenzene	10.22	180	517982	44.773	ng	96
31) Naphthalene	10.40	128	1490637	43.789	ng	100
32) Benzoic acid	9.72	122	226055	31.489	ng	97
33) 4-Chloroaniline	10.52	127	296122	19.926	ng	99
34) Hexachlorobutadiene	10.70	225	301971	44.728	ng	97
35) Caprolactam	11.29	113	157393m	42.996	ng	
36) 4-Chloro-3-methylphenol	11.66	107	519713	46.279	ng	99
37) 2-Methylnaphthalene	12.03	142	1072112	44.372	ng	98
38) 1-Methylnaphthalene	12.25	142	1009265	44.504	ng	97
40) 1,2,4,5-Tetrachlorobenzene	12.41	216	524926	45.501	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.39	237	624071	101.751	nq	99
43) 2,4,6-Trichlorophenol	12.65	196	376146	46.070	nq	98
44) 2,4,5-Trichlorophenol	12.73	196	416405	44.044	nq	99
46) 1,1'-Biphenyl	13.06	154	1326932	46.263	nq	100
47) 2-Chloronaphthalene	13.09	162	1051077	44.800	nq	99
48) 2-Nitroaniline	13.30	65	343628	45.880	nq	96
49) Acenaphthylene	13.95	152	1653471	44.155	nq	98
50) Dimethylphthalate	13.70	163	1324532	44.170	nq	99
51) 2,6-Dinitrotoluene	13.81	165	297687	45.711	nq	100
52) Acenaphthene	14.30	154	975294	44.459	nq	98
53) 3-Nitroaniline	14.14	138	185427	24.936	nq	92
54) 2,4-Dinitrophenol	14.36	184	327707	85.176	nq	98
55) Dibenzofuran	14.63	168	1542631	43.677	nq	99
56) 4-Nitrophenol	14.47	139	471523	79.861	nq	97
57) 2,4-Dinitrotoluene	14.61	165	398125	44.912	nq	97
58) Fluorene	15.29	166	1133424	40.799	nq	98
59) 2,3,4,6-Tetrachlorophenol	14.87	232	344602	45.917	nq	99
60) Diethylphthalate	15.08	149	1279780	42.593	nq	97
61) 4-Chlorophenyl-phenylether	15.29	204	574869	41.885	nq	97
62) 4-Nitroaniline	15.31	138	292853	40.987	nq	100
63) Azobenzene	15.59	77	1305817	43.660	nq	98
65) 4,6-Dinitro-2-methylphenol	15.38	198	238530	50.733	nq	98
66) n-Nitrosodiphenylamine	15.50	169	1073255	47.154	nq	100
67) 4-Bromophenyl-phenylether	16.19	248	391389	46.371	nq	96
68) Hexachlorobenzene	16.30	284	429026	47.902	nq	98
69) Atrazine	16.47	200	365557	49.233	nq	99
70) Pentachlorophenol	16.65	266	487618	91.000	nq	98
71) Phenanthrene	17.03	178	1944191	46.647	nq	99
72) Anthracene	17.12	178	1970560	47.594	nq	100
73) Carbazole	17.40	167	1776284	43.611	nq	99
74) Di-n-butylphthalate	17.98	149	2169480	45.009	nq	99
75) Fluoranthene	19.02	202	2249794	45.222	nq	99
77) Benzidine	19.20	184	872007	56.020	nq	99
78) Pyrene	19.37	202	2200244	51.564	nq	99
80) Butylbenzylphthalate	20.26	149	880498	46.420	nq	99
81) Benzo(a)anthracene	21.09	228	1828486	45.985	nq	99
82) 3,3'-Dichlorobenzidine	21.02	252	487387	32.982	nq	99
83) Chrysene	21.13	228	1757518	46.652	nq	99
84) Bis(2-ethylhexyl)phthalate	21.03	149	1223219	44.333	nq	98
85) Di-n-octyl phthalate	21.90	149	2002862	41.386	nq	99
87) Indeno(1,2,3-cd)pyrene	25.52	276	1989760	49.179	nq	99
88) Benzo(b)fluoranthene	22.64	252	1781704	51.076	nq	98
89) Benzo(k)fluoranthene	22.69	252	1666437	50.278	nq	99
90) Benzo(a)pyrene	23.20	252	1577012	48.241	nq	99
91) Dibenzo(a,h)anthracene	25.53	278	1632354	50.297	nq	98
92) Benzo(g,h,i)perylene	26.19	276	1681235	50.017	nq	98

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

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