

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122019\
 Data File : BP001348.D
 Acq On : 20 Dec 2019 17:45
 Operator : JU
 Sample : PB125625BS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB125625BS

Manual Integrations
 APPROVED

mohammad
 12/23/2019 11:56:12 AM

Quant Time: Dec 20 23:55:59 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 20 03:20:39 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.30	152	42701	20.00	ng	0.00
21) Naphthalene-d8	10.33	136	159175	20.00	ng	0.00
39) Acenaphthene-d10	13.20	164	93536	20.00	ng	0.00
64) Phenanthrene-d10	15.60	188	180523	20.00	ng	0.00
76) Chrysene-d12	19.25	240	148334	20.00	ng	0.00
87) Perylene-d12	21.02	264	153359	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	6.21	112	205232	77.68	ng	0.01
7) Phenol-d6	7.75	99	267590	77.62	ng	0.00
23) Nitrobenzene-d5	9.19	82	338221	81.60	ng	0.00
42) 2,4,6-Tribromophenol	14.50	330	82861	70.86	ng	0.00
45) 2-Fluorobiphenyl	12.12	172	593950	80.29	ng	0.00
79) Terphenyl-d14	17.89	244	728895	84.16	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.60	88	46522	42.786	ng	97
3) Pyridine	3.44	79	112185	37.769	ng	94
4) n-Nitrosodimethylamine	3.36	42	85034	45.212	ng	98
6) Aniline	7.78	93	146611	33.782	ng	# 82
8) 2-Chlorophenol	7.96	128	120056	45.199	ng	98
9) Benzaldehyde	7.59	77	94677	39.148	ng	98
10) Phenol	7.77	94	170123	43.188	ng	97
11) bis(2-Chloroethyl)ether	7.90	93	119594	42.915	ng	99
12) 1,3-Dichlorobenzene	8.20	146	139101	42.044	ng	100
13) 1,4-Dichlorobenzene	8.33	146	140276	41.616	ng	99
14) 1,2-Dichlorobenzene	8.56	146	135323	42.141	ng	97
15) Benzyl Alcohol	8.53	79	145523	44.985	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.76	45	173215	39.246	ng	99
17) 2-Methylphenol	8.73	107	104946	44.712	ng	98
18) Hexachloroethane	9.10	117	58524	40.814	ng	95
19) n-Nitroso-di-n-propylamine	8.97	70	103803	42.413	ng	99
20) 3+4-Methylphenols	8.98	107	139055	44.645	ng	96
22) Acetophenone	8.95	105	195403	38.820	ng	# 99
24) Nitrobenzene	9.22	77	162904	39.931	ng	100
25) Isophorone	9.61	82	273728	41.070	ng	100
26) 2-Nitrophenol	9.72	139	61875	42.533	ng	96
27) 2,4-Dimethylphenol	9.82	122	106630	49.836	ng	97
28) bis(2-Chloroethoxy)methane	9.98	93	148155	37.218	ng	99
29) 2,4-Dichlorophenol	10.12	162	108320	40.886	ng	99
30) 1,2,4-Trichlorobenzene	10.25	180	125763	38.592	ng	99
31) Naphthalene	10.37	128	353173	40.647	ng	100
32) Benzoic acid	10.00	122	66201	39.287	ng	97
33) 4-Chloroaniline	10.46	127	74817	21.055	ng	97
34) Hexachlorobutadiene	10.59	225	82657	36.366	ng	97
35) Caprolactam	11.04	113	32627m	41.388	ng	
36) 4-Chloro-3-methylphenol	11.28	107	126615	40.897	ng	95
37) 2-Methylnaphthalene	11.50	142	242854	40.486	ng	99
38) 1-Methylnaphthalene	11.66	142	229595	40.712	ng	96
40) 1,2,4,5-Tetrachlorobenzene	11.78	216	129898	37.559	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	11.77	237	162041	94.730	nq	98
43) 2,4,6-Trichlorophenol	11.97	196	83273	39.151	nq	100
44) 2,4,5-Trichlorophenol	12.03	196	90519	41.420	nq	98
46) 1,1'-Biphenyl	12.27	154	303502	38.774	nq	99
47) 2-Chloronaphthalene	12.29	162	236049	37.304	nq	99
48) 2-Nitroaniline	12.46	65	95888	37.843	nq	98
49) Acenaphthylene	12.97	152	374272	40.461	nq	99
50) Dimethylphthalate	12.80	163	286391	37.321	nq	99
51) 2,6-Dinitrotoluene	12.87	165	61683	39.588	nq	96
52) Acenaphthene	13.26	154	214681	38.520	nq	99
53) 3-Nitroaniline	13.13	138	39781	23.787	nq	99
54) 2,4-Dinitrophenol	13.32	184	51607	85.774	nq	96
55) Dibenzofuran	13.54	168	341040	39.055	nq	98
56) 4-Nitrophenol	13.45	139	103927	86.906	nq	91
57) 2,4-Dinitrotoluene	13.53	165	84164	38.863	nq	93
58) Fluorene	14.10	166	274074	39.241	nq	98
59) 2,3,4,6-Tetrachlorophenol	13.75	232	74739	39.835	nq	90
60) Diethylphthalate	13.96	149	290472	36.701	nq	99
61) 4-Chlorophenyl-phenylether	14.12	204	136264	37.114	nq	97
62) 4-Nitroaniline	12.46	138	74090	38.256	nq	95
63) Azobenzene	14.38	77	320353	38.447	nq	99
65) 4,6-Dinitro-2-methylphenol	14.20	198	35886	44.386	nq	97
66) n-Nitrosodiphenylamine	14.32	169	230230	40.228	nq	100
67) 4-Bromophenyl-phenylether	14.92	248	80381	37.661	nq	96
68) Hexachlorobenzene	15.00	284	93045	38.293	nq	94
69) Atrazine	15.21	200	86820	42.233	nq	96
70) Pentachlorophenol	15.32	266	109084	87.544	nq	93
71) Phenanthrene	15.64	178	408831	39.695	nq	100
72) Anthracene	15.72	178	424380	41.623	nq	99
73) Carbazole	15.96	167	366366	40.304	nq	100
74) Di-n-butylphthalate	16.52	149	454961	36.472	nq	100
75) Fluoranthene	17.35	202	461986	40.115	nq	98
77) Benzidine	17.54	184	345207	79.932	nq	100
78) Pyrene	17.66	202	464054	40.292	nq	100
80) Butylbenzylphthalate	18.54	149	198801	37.025	nq	96
81) Benzo(a)anthracene	19.23	228	423953	39.223	nq	99
82) 3,3'-Dichlorobenzidine	19.20	252	86805	23.128	nq	99
83) Chrysene	19.28	228	406654	38.978	nq	98
84) Bis(2-ethylhexyl)phthalate	19.29	149	279114	35.425	nq	99
85) Di-n-octyl phthalate	20.07	149	469338	35.940	nq	100
86) Indeno(1,2,3-cd)pyrene	22.66	276	433198	38.446	nq	99
88) Benzo(b)fluoranthene	20.53	252	398220	39.539	nq	98
89) Benzo(k)fluoranthene	20.57	252	386308	40.266	nq	99
90) Benzo(a)pyrene	20.95	252	357021	38.913	nq	98
91) Dibenzo(a,h)anthracene	22.69	278	357259	39.825	nq	99
92) Benzo(g,h,i)perylene	23.15	276	352815	41.159	nq	98

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

