

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120619\
 Data File : BP001228.D
 Acq On : 06 Dec 2019 14:22
 Operator : JU
 Sample : PB125231BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB125231BS

Manual Integrations
 APPROVED

mohammad
 12/9/2019 3:10:31 PM

Quant Time: Dec 06 14:52:14 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	100123	20.00	ng	0.00
21) Naphthalene-d8	10.34	136	412496	20.00	ng	-0.01
39) Acenaphthene-d10	13.20	164	246184	20.00	ng	0.00
64) Phenanthrene-d10	15.60	188	494524	20.00	ng	0.00
76) Chrysene-d12	19.25	240	424149	20.00	ng	-0.01
87) Perylene-d12	21.02	264	333426	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	6.21	112	748632	124.05	ng	0.00
7) Phenol-d6	7.76	99	988770	122.98	ng	0.00
23) Nitrobenzene-d5	9.19	82	568047	59.75	ng	-0.01
42) 2,4,6-Tribromophenol	14.50	330	297923	126.26	ng	0.00
45) 2-Fluorobiphenyl	12.12	172	977318	58.10	ng	-0.01
79) Terphenyl-d14	17.89	244	1306539	56.99	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.58	88	70239	24.918	ng	96
3) Pyridine	3.43	79	175328	22.415	ng	95
4) n-Nitrosodimethylamine	3.33	42	152894	45.171	ng	90
6) Aniline	7.78	93	298850	28.035	ng	# 1
8) 2-Chlorophenol	7.97	128	270651	43.347	ng	97
9) Benzaldehyde	7.59	77	52872	5.489	ng	97
10) Phenol	7.78	94	386997	42.317	ng	98
11) bis(2-Chloroethyl)ether	7.91	93	282719	39.700	ng	99
12) 1,3-Dichlorobenzene	8.21	146	264698	35.767	ng	99
13) 1,4-Dichlorobenzene	8.33	146	274693	36.846	ng	99
14) 1,2-Dichlorobenzene	8.57	146	268216	38.228	ng	98
15) Benzyl Alcohol	8.54	79	310078	46.983	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.77	45	437119	38.179	ng	99
17) 2-Methylphenol	8.73	107	244060	42.612	ng	99
18) Hexachloroethane	9.11	117	115458	37.437	ng	95
19) n-Nitroso-di-n-propylamine	8.97	70	240789	41.505	ng	99
20) 3+4-Methylphenols	8.97	107	315576	43.709	ng	99
22) Acetophenone	8.95	105	448690	37.009	ng	98
24) Nitrobenzene	9.22	77	366794	38.797	ng	98
25) Isophorone	9.62	82	650994	38.184	ng	100
26) 2-Nitrophenol	9.73	139	141193	40.772	ng	99
27) 2,4-Dimethylphenol	9.82	122	244549	44.583	ng	97
28) bis(2-Chloroethoxy)methane	9.98	93	366852	34.295	ng	99
29) 2,4-Dichlorophenol	10.12	162	246850	39.540	ng	99
30) 1,2,4-Trichlorobenzene	10.26	180	261522	35.865	ng	99
31) Naphthalene	10.37	128	777885	36.924	ng	99
32) Benzoic acid	10.02	122	172346	43.460	ng	98
33) 4-Chloroaniline	10.47	127	234499	25.650	ng	100
34) Hexachlorobutadiene	10.60	225	167651	36.549	ng	95
35) Caprolactam	11.05	113	81959m	38.091	ng	
36) 4-Chloro-3-methylphenol	11.28	107	305107	41.220	ng	99
37) 2-Methylnaphthalene	11.50	142	550027	38.262	ng	98
38) 1-Methylnaphthalene	11.66	142	513243	37.795	ng	100
40) 1,2,4,5-Tetrachlorobenzene	11.78	216	279049	35.968	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	11.77	237	225007	62.866	ng	99
43) 2,4,6-Trichlorophenol	11.97	196	192220	37.945	ng	99
44) 2,4,5-Trichlorophenol	12.03	196	204194	38.904	ng	99
46) 1,1'-Biphenyl	12.27	154	683564	35.928	ng	99
47) 2-Chloronaphthalene	12.30	162	533811	35.125	ng	99
48) 2-Nitroaniline	12.47	65	228830	38.419	ng	99
49) Acenaphthylene	12.97	152	847421	37.200	ng	99
50) Dimethylphthalate	12.80	163	693562	37.669	ng	100
51) 2,6-Dinitrotoluene	12.88	165	147028	37.303	ng	98
52) Acenaphthene	13.26	154	492150	35.915	ng	99
53) 3-Nitroaniline	13.14	138	115651	26.121	ng	97
54) 2,4-Dinitrophenol	13.32	184	133060	82.415	ng	91
55) Dibenzofuran	13.55	168	778463	37.805	ng	99
56) 4-Nitrophenol	13.45	139	254572	81.142	ng	92
57) 2,4-Dinitrotoluene	13.54	165	202164	40.921	ng	93
58) Fluorene	14.10	166	624712	37.861	ng	98
59) 2,3,4,6-Tetrachlorophenol	13.75	232	175073	40.578	ng	98
60) Diethylphthalate	13.97	149	716283	37.571	ng	99
61) 4-Chlorophenyl-phenylether	14.12	204	315741	37.578	ng	97
62) 4-Nitroaniline	12.47	138	180383	35.140	ng	99
63) Azobenzene	14.39	77	785736	38.129	ng	99
65) 4,6-Dinitro-2-methylphenol	14.20	198	94210	44.887	ng	91
66) n-Nitrosodiphenylamine	14.32	169	552320	36.758	ng	99
67) 4-Bromophenyl-phenylether	14.92	248	192587	35.871	ng	96
68) Hexachlorobenzene	15.00	284	210658	35.690	ng	97
69) Atrazine	15.22	200	189428	41.289	ng	99
70) Pentachlorophenol	15.32	266	268045	87.131	ng	99
71) Phenanthrene	15.64	178	967978	37.232	ng	99
72) Anthracene	15.72	178	975898	38.083	ng	99
73) Carbazole	15.96	167	905704	38.841	ng	100
74) Di-n-butylphthalate	16.52	149	1213624	38.710	ng	99
75) Fluoranthene	17.35	202	1118975	40.655	ng	99
77) Benzidine	17.53	184	196096	17.906	ng	99
78) Pyrene	17.66	202	1142804	34.209	ng	99
80) Butylbenzylphthalate	18.54	149	539115	34.939	ng	98
81) Benzo(a)anthracene	19.23	228	1043092	35.470	ng	99
82) 3,3'-Dichlorobenzidine	19.20	252	348544	35.876	ng	99
83) Chrysene	19.28	228	949991	36.075	ng	99
84) Bis(2-ethylhexyl)phthalate	19.29	149	770073	36.299	ng	99
85) Di-n-octyl phthalate	20.07	149	1336706	35.470	ng	100
86) Indeno(1,2,3-cd)pyrene	22.66	276	1016991	32.769	ng	99
88) Benzo(b)fluoranthene	20.53	252	965999	47.433	ng	99
89) Benzo(k)fluoranthene	20.57	252	966339	49.265	ng	99
90) Benzo(a)pyrene	20.94	252	862167	45.961	ng	98
91) Dibenzo(a,h)anthracene	22.69	278	863254	47.024	ng	99
92) Benzo(g,h,i)perylene	23.15	276	848235	46.794	ng	98

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

