

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP122619\
 Data File : BP001418.D
 Acq On : 26 Dec 2019 13:27
 Operator : JU
 Sample : PB125701BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB125701BS

Manual Integrations
 APPROVED

mohammad
 12/27/2019 3:18:02 PM

Quant Time: Dec 27 07:42:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP121919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 24 16:04:49 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.29	152	35390	20.00	ng	0.00
21) Naphthalene-d8	10.32	136	135788	20.00	ng	0.00
39) Acenaphthene-d10	13.19	164	83844	20.00	ng	0.00
64) Phenanthrene-d10	15.58	188	165462	20.00	ng	0.00
76) Chrysene-d12	19.23	240	105971	20.00	ng	0.00
87) Perylene-d12	20.99	264	79095	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	6.22	112	273199	124.76	ng	0.02
7) Phenol-d6	7.76	99	357676	125.18	ng	0.00
23) Nitrobenzene-d5	9.18	82	261424	73.93	ng	0.00
42) 2,4,6-Tribromophenol	14.49	330	122210	116.59	ng	0.00
45) 2-Fluorobiphenyl	12.10	172	526441	79.39	ng	0.00
79) Terphenyl-d14	17.87	244	575861	93.07	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.61	88	33215	36.858	ng	95
3) Pyridine	3.45	79	78412	31.852	ng	98
4) n-Nitrosodimethylamine	3.36	42	69099	44.329	ng	100
6) Aniline	7.76	93	112124	31.173	ng	# 94
8) 2-Chlorophenol	7.96	128	103499	47.015	ng	97
9) Benzaldehyde	7.58	77	46138	23.019	ng	96
10) Phenol	7.78	94	143270	43.884	ng	84
11) bis(2-Chloroethyl)ether	7.89	93	96488	41.776	ng	98
12) 1,3-Dichlorobenzene	8.19	146	117847	42.978	ng	99
13) 1,4-Dichlorobenzene	8.32	146	119566	42.799	ng	99
14) 1,2-Dichlorobenzene	8.55	146	115441	43.376	ng	97
15) Benzyl Alcohol	8.53	79	117145	43.694	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.75	45	133122	36.393	ng	93
17) 2-Methylphenol	8.72	107	89270	45.890	ng	97
18) Hexachloroethane	9.09	117	48379	40.709	ng	97
19) n-Nitroso-di-n-propylamine	8.96	70	84703	41.758	ng	98
20) 3+4-Methylphenols	8.97	107	118612	45.948	ng	98
22) Acetophenone	8.94	105	167079	38.910	ng	# 99
24) Nitrobenzene	9.20	77	135803	39.021	ng	99
25) Isophorone	9.60	82	225530	39.666	ng	99
26) 2-Nitrophenol	9.71	139	54923	44.257	ng	98
27) 2,4-Dimethylphenol	9.82	122	92335	50.588	ng	96
28) bis(2-Chloroethoxy)methane	9.96	93	123774	36.449	ng	98
29) 2,4-Dichlorophenol	10.11	162	98646	43.648	ng	99
30) 1,2,4-Trichlorobenzene	10.24	180	112476	40.459	ng	97
31) Naphthalene	10.36	128	311474	42.022	ng	99
32) Benzoic acid	10.03	122	49262	34.911	ng	94
33) 4-Chloroaniline	10.45	127	67583	22.295	ng	99
34) Hexachlorobutadiene	10.57	225	75181	38.773	ng	92
35) Caprolactam	11.04	113	28225m	41.970	ng	
36) 4-Chloro-3-methylphenol	11.28	107	110416	41.807	ng	99
37) 2-Methylnaphthalene	11.49	142	223336	43.644	ng	100
38) 1-Methylnaphthalene	11.65	142	211598	43.983	ng	98
40) 1,2,4,5-Tetrachlorobenzene	11.76	216	123068	39.698	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	11.76	237	108479	70.748	nq	95
43) 2,4,6-Trichlorophenol	11.96	196	78395	41.118	nq	99
44) 2,4,5-Trichlorophenol	12.03	196	85307	43.548	nq	98
46) 1,1'-Biphenyl	12.26	154	284124	40.495	nq	99
47) 2-Chloronaphthalene	12.28	162	220609	38.894	nq	98
48) 2-Nitroaniline	12.45	65	83347	36.696	nq	97
49) Acenaphthylene	12.95	152	345914	41.718	nq	99
50) Dimethylphthalate	12.79	163	266084	38.683	nq	99
51) 2,6-Dinitrotoluene	12.86	165	57529	41.190	nq	85
52) Acenaphthene	13.24	154	200610	40.156	nq	99
53) 3-Nitroaniline	13.13	138	35585	23.738	nq	97
54) 2,4-Dinitrophenol	13.31	184	48582	90.080	nq	86
55) Dibenzofuran	13.53	168	327875	41.888	nq	99
56) 4-Nitrophenol	13.45	139	79494	74.159	nq	95
57) 2,4-Dinitrotoluene	13.52	165	78502	40.439	nq	# 90
58) Fluorene	14.09	166	257586	41.144	nq	99
59) 2,3,4,6-Tetrachlorophenol	13.74	232	69500	41.324	nq	91
60) Diethylphthalate	13.95	149	270499	38.128	nq	98
61) 4-Chlorophenyl-phenylether	14.10	204	136521	41.483	nq	99
62) 4-Nitroaniline	12.45	138	67009	38.600	nq	90
63) Azobenzene	14.36	77	284692	38.116	nq	95
65) 4,6-Dinitro-2-methylphenol	14.19	198	36395	49.114	nq	92
66) n-Nitrosodiphenylamine	14.30	169	221382	42.203	nq	99
67) 4-Bromophenyl-phenylether	14.90	248	80798	41.302	nq	95
68) Hexachlorobenzene	14.99	284	89075	39.996	nq	96
69) Atrazine	15.20	200	76327	40.509	nq	97
70) Pentachlorophenol	15.30	266	87689	76.779	nq	97
71) Phenanthrene	15.62	178	378675	40.113	nq	100
72) Anthracene	15.70	178	385060	41.204	nq	99
73) Carbazole	15.95	167	316751	38.018	nq	98
74) Di-n-butylphthalate	16.50	149	410985	35.946	nq	99
75) Fluoranthene	17.33	202	392102	37.146	nq	99
77) Benzidine	17.52	184	89851	29.122	nq	99
78) Pyrene	17.64	202	382631	46.504	nq	98
80) Butylbenzylphthalate	18.52	149	146243	38.124	nq	91
81) Benzo(a)anthracene	19.22	228	304884	39.483	nq	100
82) 3,3'-Dichlorobenzidine	19.19	252	94575	35.271	nq	98
83) Chrysene	19.26	228	293048	39.317	nq	99
84) Bis(2-ethylhexyl)phthalate	19.26	149	205699	36.544	nq	100
85) Di-n-octyl phthalate	20.05	149	310437	33.275	nq	97
86) Indeno(1,2,3-cd)pyrene	22.63	276	289709	35.990	nq	97
88) Benzo(b)fluoranthene	20.52	252	265012	51.019	nq	99
89) Benzo(k)fluoranthene	20.55	252	268158	54.194	nq	99
90) Benzo(a)pyrene	20.93	252	236500	49.979	nq	98
91) Dibenzo(a,h)anthracene	22.65	278	248387	53.686	nq	99
92) Benzo(g,h,i)perylene	23.11	276	239466	54.165	nq	99

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

