

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP121019\
 Data File : BP001295.D
 Acq On : 10 Dec 2019 12:30
 Operator : JU
 Sample : PB125342BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PB125342BS

Manual Integrations
 APPROVED

mohammad
 12/11/2019 12:28:40 PM

Quant Time: Dec 11 04:12:11 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	96560	20.00	ng	0.00
21) Naphthalene-d8	10.35	136	413841	20.00	ng	0.00
39) Acenaphthene-d10	13.21	164	246804	20.00	ng	0.00
64) Phenanthrene-d10	15.61	188	457863	20.00	ng	0.00
76) Chrysene-d12	19.26	240	350043	20.00	ng	0.00
87) Perylene-d12	21.03	264	303254	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	6.23	112	868557	149.23	ng	0.02
7) Phenol-d6	7.78	99	1173606	151.36	ng	0.01
23) Nitrobenzene-d5	9.20	82	677482	71.03	ng	0.00
42) 2,4,6-Tribromophenol	14.52	330	358326	151.47	ng	0.01
45) 2-Fluorobiphenyl	12.13	172	1241265	73.61	ng	0.00
79) Terphenyl-d14	17.89	244	1405613	74.29	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.60	88	80967	29.784	ng	95
3) Pyridine	3.45	79	209594	27.784	ng	95
4) n-Nitrosodimethylamine	3.36	42	184649	56.566	ng	80
6) Aniline	7.79	93	384005	37.353	ng	# 13
8) 2-Chlorophenol	7.98	128	320400	53.208	ng	96
9) Benzaldehyde	7.60	77	122871	24.296	ng	97
10) Phenol	7.79	94	456022	51.705	ng	99
11) bis(2-Chloroethyl)ether	7.92	93	316644	46.105	ng	99
12) 1,3-Dichlorobenzene	8.22	146	315301	44.176	ng	98
13) 1,4-Dichlorobenzene	8.33	146	322675	44.879	ng	98
14) 1,2-Dichlorobenzene	8.58	146	317757	46.960	ng	99
15) Benzyl Alcohol	8.55	79	384075	60.343	ng	98
16) 2,2'-oxybis(1-Chloropropan	8.77	45	491356	44.500	ng	99
17) 2-Methylphenol	8.74	107	285471	51.681	ng	96
18) Hexachloroethane	9.11	117	137768	46.320	ng	97
19) n-Nitroso-di-n-propylamine	8.99	70	293798	52.511	ng	98
20) 3+4-Methylphenols	8.99	107	381718	54.821	ng	97
22) Acetophenone	8.96	105	535258	44.006	ng	# 97
24) Nitrobenzene	9.23	77	429886	45.322	ng	98
25) Isophorone	9.62	82	770817	45.065	ng	99
26) 2-Nitrophenol	9.73	139	167003	48.069	ng	94
27) 2,4-Dimethylphenol	9.83	122	293294	53.295	ng	97
28) bis(2-Chloroethoxy)methane	9.99	93	432950	40.343	ng	99
29) 2,4-Dichlorophenol	10.13	162	302467	48.291	ng	98
30) 1,2,4-Trichlorobenzene	10.26	180	321149	43.899	ng	97
31) Naphthalene	10.38	128	922999	43.670	ng	99
32) Benzoic acid	10.06	122	199766	50.211	ng	95
33) 4-Chloroaniline	10.48	127	256951	28.015	ng	98
34) Hexachlorobutadiene	10.60	225	213320	46.353	ng	97
35) Caprolactam	11.08	113	86848m	40.232	ng	
36) 4-Chloro-3-methylphenol	11.30	107	357751	48.175	ng	96
37) 2-Methylnaphthalene	11.52	142	672380	46.621	ng	99
38) 1-Methylnaphthalene	11.68	142	624850	45.865	ng	100
40) 1,2,4,5-Tetrachlorobenzene	11.79	216	361289	46.451	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	11.78	237	410827	114.495	ng	97
43) 2,4,6-Trichlorophenol	11.98	196	239069	47.075	ng	98
44) 2,4,5-Trichlorophenol	12.05	196	253214	48.123	ng	98
46) 1,1'-Biphenyl	12.29	154	856276	44.893	ng	98
47) 2-Chloronaphthalene	12.30	162	654386	42.950	ng	99
48) 2-Nitroaniline	12.47	65	260353	43.601	ng	97
49) Acenaphthylene	12.98	152	1014191	44.409	ng	100
50) Dimethylphthalate	12.82	163	813427	44.068	ng	99
51) 2,6-Dinitrotoluene	12.89	165	172314	43.609	ng	96
52) Acenaphthene	13.27	154	591265	43.040	ng	100
53) 3-Nitroaniline	13.15	138	120514	27.151	ng	99
54) 2,4-Dinitrophenol	13.33	184	154539	94.351	ng	86
55) Dibenzofuran	13.56	168	935243	45.305	ng	98
56) 4-Nitrophenol	13.47	139	263973	83.927	ng	89
57) 2,4-Dinitrotoluene	13.55	165	229880	46.414	ng	# 86
58) Fluorene	14.12	166	751225	45.414	ng	100
59) 2,3,4,6-Tetrachlorophenol	13.76	232	207523	47.978	ng	99
60) Diethylphthalate	13.98	149	837078	43.797	ng	99
61) 4-Chlorophenyl-phenylether	14.13	204	388227	46.089	ng	98
62) 4-Nitroaniline	12.47	138	202805	39.409	ng	97
63) Azobenzene	14.39	77	901222	43.623	ng	95
65) 4,6-Dinitro-2-methylphenol	14.22	198	105816	53.116	ng	92
66) n-Nitrosodiphenylamine	14.33	169	635794	45.701	ng	99
67) 4-Bromophenyl-phenylether	14.93	248	230022	46.274	ng	100
68) Hexachlorobenzene	15.02	284	251325	45.989	ng	97
69) Atrazine	15.23	200	221499	52.145	ng	98
70) Pentachlorophenol	15.33	266	302622	106.247	ng	99
71) Phenanthrene	15.65	178	1090246	45.292	ng	99
72) Anthracene	15.73	178	1104873	46.568	ng	99
73) Carbazole	15.97	167	957678	44.359	ng	99
74) Di-n-butylphthalate	16.52	149	1336816	46.053	ng	99
75) Fluoranthene	17.36	202	1180400	46.321	ng	99
77) Benzidine	17.55	184	360511	39.889	ng	98
78) Pyrene	17.66	202	1176800	42.684	ng	99
80) Butylbenzylphthalate	18.55	149	544168	42.733	ng	97
81) Benzo(a)anthracene	19.25	228	1024095	42.196	ng	100
82) 3,3'-Dichlorobenzidine	19.21	252	348771	43.500	ng	99
83) Chrysene	19.29	228	970416	44.652	ng	99
84) Bis(2-ethylhexyl)phthalate	19.29	149	805524	46.009	ng	99
85) Di-n-octyl phthalate	20.07	149	1329520	42.748	ng	100
86) Indeno(1,2,3-cd)pyrene	22.67	276	968798	37.825	ng	98
88) Benzo(b)fluoranthene	20.55	252	919231	49.627	ng	98
89) Benzo(k)fluoranthene	20.58	252	906678	50.822	ng	99
90) Benzo(a)pyrene	20.96	252	816273	47.844	ng	99
91) Dibenzo(a,h)anthracene	22.70	278	823425	49.318	ng	98
92) Benzo(g,h,i)perylene	23.17	276	795678	48.262	ng	97

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

