

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021219\
 Data File : BM018773.D
 Acq On : 12 Feb 2019 15:22
 Operator : JU/SJ
 Sample : PB117038BSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB117038BSD

Manual Integrations
 APPROVED

Sohil
 2/13/2019 12:20:43 PM

Quant Time: Feb 13 02:41:47 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 11 16:02:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.72	152	335186	20.00	ng	0.00
21) Naphthalene-d8	10.50	136	1378617	20.00	ng	0.00
39) Acenaphthene-d10	14.37	164	791910	20.00	ng	0.00
64) Phenanthrene-d10	17.12	188	1837829	20.00	ng	0.00
76) Chrysene-d12	21.31	240	2078547	20.00	ng	0.00
87) Perylene-d12	23.57	264	2041833	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.30	112	2923263	142.90	ng	0.00
7) Phenol-d6	6.89	99	4145813	143.86	ng	0.00
23) Nitrobenzene-d5	8.89	82	2142537	71.86	ng	0.00
42) 2,4,6-Tribromophenol	15.87	330	1788832	156.71	ng	0.00
45) 2-Fluorobiphenyl	12.98	172	4232378	70.95	ng	0.00
79) Terphenyl-d14	19.75	244	7898010	78.39	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.26	88	364797	40.942	ng	99
3) Pyridine	3.66	79	1111562	45.736	ng	98
4) n-Nitrosodimethylamine	3.58	42	302618	48.945	ng	100
6) Aniline	7.06	93	1585140	44.891	ng	99
8) 2-Chlorophenol	7.28	128	959076	42.701	ng	97
9) Benzaldehyde	6.87	77	859466	63.716	ng	99
10) Phenol	6.92	94	1311574	43.253	ng	99
11) bis(2-Chloroethyl)ether	7.16	93	973150	42.071	ng	98
12) 1,3-Dichlorobenzene	7.60	146	1040613	41.204	ng	99
13) 1,4-Dichlorobenzene	7.75	146	1055993	41.071	ng	99
14) 1,2-Dichlorobenzene	8.06	146	1020427	41.318	ng	99
15) Benzyl Alcohol	7.96	79	988456	43.165	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.23	45	915538	40.894	ng	98
17) 2-Methylphenol	8.16	107	831904	43.104	ng	98
18) Hexachloroethane	8.78	117	395492	42.138	ng	99
19) n-Nitroso-di-n-propylamine	8.53	70	937605	42.019	ng	100
20) 3+4-Methylphenols	8.49	107	1150617	43.174	ng	98
22) Acetophenone	8.55	105	1554540	41.061	ng	# 99
24) Nitrobenzene	8.93	77	1312097	42.396	ng	98
25) Isophorone	9.45	82	2360692	49.622	ng	100
26) 2-Nitrophenol	9.63	139	541284	43.918	ng	99
27) 2,4-Dimethylphenol	9.69	122	875157	48.617	ng	98
28) bis(2-Chloroethoxy)methane	9.93	93	1360921	44.238	ng	99
29) 2,4-Dichlorophenol	10.16	162	905136	43.791	ng	99
30) 1,2,4-Trichlorobenzene	10.36	180	991828	41.551	ng	99
31) Naphthalene	10.56	128	3019495	44.909	ng	99
32) Benzoic acid	9.85	122	615818	39.218	ng	98
33) 4-Chloroaniline	10.68	127	1321854	43.980	ng	99
34) Hexachlorobutadiene	10.82	225	550037	39.699	ng	99
35) Caprolactam	11.50	113	323168m	38.310	ng	
36) 4-Chloro-3-methylphenol	11.80	107	1075341	43.151	ng	99
37) 2-Methylnaphthalene	12.17	142	2209039	46.665	ng	100
38) 1-Methylnaphthalene	12.39	142	2094276	45.241	ng	100
40) 1,2,4,5-Tetrachlorobenzene	12.54	216	1057866	41.759	ng	99

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM021219\
 Data File : BM018773.D
 Acq On : 12 Feb 2019 15:22
 Operator : JU/SJ
 Sample : PB117038BSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 PB117038BSD

Manual Integrations
 APPROVED

Sohil
 2/13/2019 12:20:43 PM

Quant Time: Feb 13 02:41:47 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 11 16:02:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	12.51	237	1436899	110.942	ng	100
43) 2,4,6-Trichlorophenol	12.79	196	741624	44.223	ng	99
44) 2,4,5-Trichlorophenol	12.86	196	792013	45.059	ng	97
46) 1,1'-Biphenyl	13.19	154	2769621	40.631	ng	100
47) 2-Chloronaphthalene	13.24	162	2170314	41.221	ng	100
48) 2-Nitroaniline	13.46	65	783844	44.826	ng	97
49) Acenaphthylene	14.09	152	3566859	45.497	ng	100
50) Dimethylphthalate	13.83	163	2845745	43.119	ng	100
51) 2,6-Dinitrotoluene	13.96	165	629935	44.062	ng	99
52) Acenaphthene	14.43	154	2043168	42.503	ng	99
53) 3-Nitroaniline	14.29	138	676740	41.894	ng	97
54) 2,4-Dinitrophenol	14.50	184	816161	89.251	ng	96
55) Dibenzofuran	14.77	168	3339968	42.271	ng	99
56) 4-Nitrophenol	14.61	139	1143916	88.710	ng	99
57) 2,4-Dinitrotoluene	14.75	165	902164	45.800	ng	98
58) Fluorene	15.42	166	2749574	45.487	ng	99
59) 2,3,4,6-Tetrachlorophenol	14.99	232	733098	47.185	ng	100
60) Diethylphthalate	15.20	149	2952860	43.374	ng	99
61) 4-Chlorophenyl-phenylether	15.42	204	1374609	40.559	ng	97
62) 4-Nitroaniline	15.47	138	693029	43.762	ng	98
63) Azobenzene	15.71	77	3128839	42.864	ng	99
65) 4,6-Dinitro-2-methylphenol	15.51	198	522593	40.500	ng	98
66) n-Nitrosodiphenylamine	15.63	169	2419825	41.675	ng	99
67) 4-Bromophenyl-phenylether	16.31	248	831168	40.404	ng	99
68) Hexachlorobenzene	16.42	284	956250	39.992	ng	97
69) Atrazine	16.59	200	827330	44.197	ng	98
70) Pentachlorophenol	16.76	266	1272104	82.627	ng	100
71) Phenanthrene	17.16	178	4362819	44.173	ng	100
72) Anthracene	17.25	178	4456789	46.360	ng	100
73) Carbazole	17.53	167	4117380	40.995	ng	99
74) Di-n-butylphthalate	18.08	149	5121124	44.335	ng	99
75) Fluoranthene	19.18	202	5215078	45.711	ng	98
77) Benzidine	19.39	184	6291564	134.401	ng	99
78) Pyrene	19.54	202	5340649	44.240	ng	99
80) Butylbenzylphthalate	20.45	149	2540892	46.184	ng	97
81) Benzo(a)anthracene	21.30	228	5468983	45.138	ng	100
82) 3,3'-Dichlorobenzidine	21.24	252	2169999	45.317	ng	100
83) Chrysene	21.35	228	5065855	43.621	ng	100
84) Bis(2-ethylhexyl)phthalate	21.22	149	3657016	45.416	ng	100
85) Di-n-octyl phthalate	22.10	149	6378069	47.128	ng	100
86) Indeno(1,2,3-cd)pyrene	25.87	276	5643283	42.089	ng	100
88) Benzo(b)fluoranthene	22.89	252	5571333	47.691	ng	99
89) Benzo(k)fluoranthene	22.94	252	5219753	44.881	ng	100
90) Benzo(a)pyrene	23.47	252	5173765	46.751	ng	99
91) Dibenzo(a,h)anthracene	25.88	278	4822202	43.981	ng	98
92) Benzo(g,h,i)perylene	26.57	276	4412614	43.229	ng	99

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

