

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080719\
 Data File : BF116041.D
 Acq On : 7 Aug 2019 10:58
 Operator : HP/JU
 Sample : PB122007BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB122007BS

Manual Integrations
 APPROVED

mohammad
 8/8/2019 4:46:41 PM

Quant Time: Aug 07 13:09:37 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 11:13:18 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	139187	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	569291	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	309268	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	543634	20.00	ng	0.00
76) Chrysene-d12	14.01	240	386117	20.00	ng	0.00
87) Perylene-d12	15.47	264	315473	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.48	112	874514	105.18	ng	0.01
7) Phenol-d6	6.49	99	1115400	106.33	ng	0.00
23) Nitrobenzene-d5	7.41	82	721494	73.16	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	320712	106.96	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	1270724	76.96	ng	0.00
79) Terphenyl-d14	12.94	244	1465296	79.97	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.69	88	161256	38.391	ng	98
3) Pyridine	3.45	79	333878	28.456	ng	98
4) n-Nitrosodimethylamine	3.39	42	161182	34.810	ng	99
6) Aniline	6.51	93	416788	27.743	ng	99
8) 2-Chlorophenol	6.63	128	361955	39.369	ng	97
9) Benzaldehyde	6.40	77	120151	16.600	ng	98
10) Phenol	6.50	94	461951	37.396	ng	88
11) bis(2-Chloroethyl)ether	6.58	93	346056	38.103	ng	99
12) 1,3-Dichlorobenzene	6.79	146	379736	35.738	ng	99
13) 1,4-Dichlorobenzene	6.86	146	386369	36.317	ng	98
14) 1,2-Dichlorobenzene	7.02	146	358494	36.595	ng	99
15) Benzyl Alcohol	6.99	79	309596	38.890	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.12	45	461644	37.266	ng	97
17) 2-Methylphenol	7.10	107	306793	39.408	ng	99
18) Hexachloroethane	7.36	117	140017	35.746	ng	95
19) n-Nitroso-di-n-propylamine	7.26	70	250562	38.546	ng	97
20) 3+4-Methylphenols	7.26	107	373678	40.562	ng	96
22) Acetophenone	7.25	105	498437	39.922	ng	99
24) Nitrobenzene	7.43	77	405080	38.039	ng	99
25) Isophorone	7.67	82	730367	38.729	ng	99
26) 2-Nitrophenol	7.75	139	199849	39.812	ng	95
27) 2,4-Dimethylphenol	7.78	122	327528	43.368	ng	100
28) bis(2-Chloroethoxy)methane	7.87	93	450303	38.524	ng	100
29) 2,4-Dichlorophenol	7.99	162	317831	39.858	ng	98
30) 1,2,4-Trichlorobenzene	8.07	180	335336	36.892	ng	99
31) Naphthalene	8.15	128	1041437	38.875	ng	99
32) Benzoic acid	7.93	122	192983	36.244	ng	99
33) 4-Chloroaniline	8.20	127	330479	27.609	ng	97
34) Hexachlorobutadiene	8.26	225	186673	35.654	ng	97
35) Caprolactam	8.57	113	99791m	38.693	ng	
36) 4-Chloro-3-methylphenol	8.69	107	334787	40.357	ng	95
37) 2-Methylnaphthalene	8.84	142	699027	39.620	ng	100
38) 1-Methylnaphthalene	8.94	142	660557	39.575	ng	100
40) 1,2,4,5-Tetrachlorobenzene	9.01	216	323020	39.069	ng	100

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.99	237	148949	43.081	nq	100
43) 2,4,6-Trichlorophenol	9.12	196	208386	36.617	nq	99
44) 2,4,5-Trichlorophenol	9.17	196	228493	38.841	nq	97
46) 1,1'-Biphenyl	9.30	154	864247	39.582	nq	98
47) 2-Chloronaphthalene	9.33	162	678336	37.328	nq	99
48) 2-Nitroaniline	9.43	65	220960	36.786	nq	98
49) Acenaphthylene	9.75	152	1025139	38.667	nq	99
50) Dimethylphthalate	9.60	163	821079	38.992	nq	99
51) 2,6-Dinitrotoluene	9.67	165	181660	39.697	nq	99
52) Acenaphthene	9.92	154	617597	37.972	nq	99
53) 3-Nitroaniline	9.84	138	154981	27.409	nq	94
54) 2,4-Dinitrophenol	9.95	184	149097	65.615	nq	92
55) Dibenzofuran	10.09	168	916883	38.764	nq	99
56) 4-Nitrophenol	10.01	139	251727	69.494	nq	96
57) 2,4-Dinitrotoluene	10.07	165	237517	38.983	nq	99
58) Fluorene	10.43	166	725527	39.868	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.21	232	188911	38.170	nq	97
60) Diethylphthalate	10.30	149	788381	40.101	nq	98
61) 4-Chlorophenyl-phenylether	10.42	204	364637	39.564	nq	98
62) 4-Nitroaniline	10.46	138	199212	34.091	nq	99
63) Azobenzene	10.58	77	811425	40.127	nq	97
65) 4,6-Dinitro-2-methylphenol	10.49	198	116025	40.274	nq	98
66) n-Nitrosodiphenylamine	10.54	169	653873	39.961	nq	100
67) 4-Bromophenyl-phenylether	10.90	248	225805	38.801	nq	96
68) Hexachlorobenzene	10.98	284	252022	37.451	nq	98
69) Atrazine	11.06	200	204180	37.930	nq	99
70) Pentachlorophenol	11.18	266	228289	69.875	nq	99
71) Phenanthrene	11.39	178	1073852	38.639	nq	99
72) Anthracene	11.45	178	1106139	39.327	nq	99
73) Carbazole	11.60	167	1020881	40.007	nq	99
74) Di-n-butylphthalate	11.92	149	1249360	41.971	nq	99
75) Fluoranthene	12.58	202	1152480	39.611	nq	99
77) Benzidine	12.70	184	358407	27.475	nq	97
78) Pyrene	12.81	202	1151402	39.559	nq	99
80) Butylbenzylphthalate	13.42	149	532869	43.280	nq	94
81) Benzo(a)anthracene	14.00	228	962348	38.994	nq	100
82) 3,3'-Dichlorobenzidine	13.96	252	314249	36.651	nq	99
83) Chrysene	14.03	228	938779	39.182	nq	99
84) Bis(2-ethylhexyl)phthalate	13.97	149	756965	47.312	nq	99
85) Di-n-octyl phthalate	14.59	149	1191818	46.898	nq	99
86) Indeno(1,2,3-cd)pyrene	16.95	276	765668	38.048	nq	100
88) Benzo(b)fluoranthene	15.05	252	846367	46.399	nq	100
89) Benzo(k)fluoranthene	15.08	252	804432	45.277	nq	100
90) Benzo(a)pyrene	15.42	252	764000	46.854	nq	100
91) Dibenzo(a,h)anthracene	16.96	278	646191	45.782	nq	100
92) Benzo(g,h,i)perylene	17.39	276	620206	44.742	nq	99

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

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