

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101518\
 Data File : BF109888.D
 Acq On : 15 Oct 2018 19:26
 Operator : JU/SJ
 Sample : PB113836BSD
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB113836BSD

Manual Integrations
 APPROVED

Sohil
 10/16/2018 9:57:31 AM

Quant Time: Oct 16 01:41:59 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 15 18:09:48 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	243248	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	945797	20.00	ng	0.00
39) Acenaphthene-d10	10.02	164	433318	20.00	ng	0.00
64) Phenanthrene-d10	11.51	188	770509	20.00	ng	0.00
76) Chrysene-d12	14.15	240	456310	20.00	ng	0.00
87) Perylene-d12	15.67	264	511213	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.60	112	1603820	104.51	ng	0.01
7) Phenol-d6	6.60	99	2029502	106.51	ng	0.00
23) Nitrobenzene-d5	7.54	82	1235617	88.01	ng	0.00
42) 2,4,6-Tribromophenol	10.81	330	437143	116.52	ng	0.00
45) 2-Fluorobiphenyl	9.33	172	2167076	79.80	ng	0.00
79) Terphenyl-d14	13.09	244	1826016	84.03	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.91	88	265055	30.411	ng	92
3) Pyridine	3.66	79	752382	38.720	ng	90
4) n-Nitrosodimethylamine	3.62	42	365390	46.096	ng	# 99
6) Aniline	6.64	93	536023	20.874	ng	# 52
8) 2-Chlorophenol	6.76	128	721603	41.980	ng	98
9) Benzaldehyde	6.53	77	357623	28.880	ng	100
10) Phenol	6.61	94	839555	40.808	ng	# 62
11) bis(2-Chloroethyl)ether	6.71	93	697268	40.534	ng	92
12) 1,3-Dichlorobenzene	6.92	146	757675	40.198	ng	95
13) 1,4-Dichlorobenzene	6.99	146	764409	40.269	ng	98
14) 1,2-Dichlorobenzene	7.15	146	708443	40.528	ng	99
15) Benzyl Alcohol	7.11	79	619832	42.281	ng	92
16) 2,2'-oxybis(1-Chloropropan	7.25	45	1165508	41.280	ng	67
17) 2-Methylphenol	7.22	107	598604	43.186	ng	# 81
18) Hexachloroethane	7.49	117	280747	41.851	ng	96
19) n-Nitroso-di-n-propylamine	7.39	70	489454	40.005	ng	96
20) 3+4-Methylphenols	7.37	107	737692	41.940	ng	# 89
22) Acetophenone	7.39	105	968072	44.136	ng	98
24) Nitrobenzene	7.56	77	729462	47.826	ng	96
25) Isophorone	7.80	82	1365547	49.499	ng	95
26) 2-Nitrophenol	7.87	139	339369	53.664	ng	92
27) 2,4-Dimethylphenol	7.90	122	677021	50.950	ng	95
28) bis(2-Chloroethoxy)methane	8.00	93	877550	46.350	ng	99
29) 2,4-Dichlorophenol	8.11	162	588490	45.653	ng	100
30) 1,2,4-Trichlorobenzene	8.20	180	625364	43.328	ng	96
31) Naphthalene	8.28	128	2019392	46.561	ng	99
32) Benzoic acid	7.99	122	210918	25.475	ng	93
33) 4-Chloroaniline	8.32	127	196677	10.088	ng	99
34) Hexachlorobutadiene	8.40	225	336577	42.289	ng	98
35) Caprolactam	8.69	113	196192m	42.855	ng	
36) 4-Chloro-3-methylphenol	8.80	107	617937	45.441	ng	97
37) 2-Methylnaphthalene	8.97	142	1380096	49.142	ng	98
38) 1-Methylnaphthalene	9.07	142	1313932	48.283	ng	98
40) 1,2,4,5-Tetrachlorobenzene	9.14	216	561109	42.041	ng	# 98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.13	237	679607	105.498	nq	99
43) 2,4,6-Trichlorophenol	9.25	196	394202	47.235	nq	96
44) 2,4,5-Trichlorophenol	9.29	196	408553	47.168	nq	94
46) 1,1'-Biphenyl	9.44	154	1610105	41.705	nq	97
47) 2-Chloronaphthalene	9.46	162	1238072	43.536	nq	97
48) 2-Nitroaniline	9.56	65	374404	51.303	nq	97
49) Acenaphthylene	9.89	152	1941400	46.993	nq	97
50) Dimethylphthalate	9.73	163	1401921	45.691	nq	98
51) 2,6-Dinitrotoluene	9.80	165	305430	52.410	nq	95
52) Acenaphthene	10.06	154	1247329	47.424	nq	99
53) 3-Nitroaniline	9.96	138	128133	18.075	nq	98
54) 2,4-Dinitrophenol	10.07	184	150799	81.091	nq	# 79
55) Dibenzofuran	10.23	168	1685971	44.947	nq	97
56) 4-Nitrophenol	10.12	139	505960	93.538	nq	87
57) 2,4-Dinitrotoluene	10.20	165	385530	51.953	nq	# 89
58) Fluorene	10.57	166	1266426	46.380	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.34	232	327937	47.772	nq	95
60) Diethylphthalate	10.44	149	1343612	44.337	nq	97
61) 4-Chlorophenyl-phenylether	10.56	204	599114	42.121	nq	97
62) 4-Nitroaniline	10.59	138	297570	44.344	nq	92
63) Azobenzene	10.72	77	1379224	43.565	nq	95
65) 4,6-Dinitro-2-methylphenol	10.61	198	117830	43.165	nq	98
66) n-Nitrosodiphenylamine	10.68	169	1139391	44.399	nq	96
67) 4-Bromophenyl-phenylether	11.05	248	371280	44.463	nq	93
68) Hexachlorobenzene	11.12	284	390026	43.801	nq	# 92
69) Atrazine	11.20	200	370920	46.272	nq	98
70) Pentachlorophenol	11.31	266	347967	68.665	nq	96
71) Phenanthrene	11.54	178	1821853	45.672	nq	99
72) Anthracene	11.59	178	1911144	47.655	nq	99
73) Carbazole	11.74	167	1627127	41.283	nq	99
74) Di-n-butylphthalate	12.06	149	1904483	43.297	nq	99
75) Fluoranthene	12.73	202	1744943	43.789	nq	99
77) Benzidine	12.84	184	430575	24.611	nq	97
78) Pyrene	12.96	202	1727010	51.530	nq	100
80) Butylbenzylphthalate	13.56	149	674967	49.815	nq	98
81) Benzo(a)anthracene	14.14	228	1304577	48.537	nq	99
82) 3,3'-Dichlorobenzidine	14.10	252	203452	21.591	nq	98
83) Chrysene	14.18	228	1205544	47.122	nq	99
84) Bis(2-ethylhexyl)phthalate	14.12	149	840074	47.118	nq	96
85) Di-n-octyl phthalate	14.74	149	1325799	52.745	nq	98
86) Indeno(1,2,3-cd)pyrene	17.23	276	1747492	84.835	nq	# 93
88) Benzo(b)fluoranthene	15.22	252	1284338	43.577	nq	97
89) Benzo(k)fluoranthene	15.25	252	1218305	41.938	nq	# 95
90) Benzo(a)pyrene	15.61	252	1304027	48.108	nq	97
91) Dibenzo(a,h)anthracene	17.26	278	1428812	59.478	nq	# 96
92) Benzo(g,h,i)perylene	17.72	276	1492325	61.485	nq	# 94

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

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