

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102518\
 Data File : BF110166.D
 Acq On : 25 Oct 2018 17:54
 Operator : JU/SJ
 Sample : J5636-02MS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HP-GRID-BMS

Manual Integrations
 APPROVED

Sohil
 10/26/2018 3:49:18 PM

Quant Time: Oct 26 04:22:55 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 23 15:06:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	156450	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	629390	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	299850	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	543574	20.00	ng	-0.01
76) Chrysene-d12	14.15	240	296489	20.00	ng	-0.01
87) Perylene-d12	15.67	264	275136	20.00	ng	-0.02

System Monitoring Compounds

5) 2-Fluorophenol	5.60	112	1298181	135.12	ng	0.00
7) Phenol-d6	6.60	99	1665530	135.54	ng	0.00
23) Nitrobenzene-d5	7.53	82	1051339	99.08	ng	-0.02
42) 2,4,6-Tribromophenol	10.81	330	391453	131.79	ng	0.00
45) 2-Fluorobiphenyl	9.33	172	1753857	91.85	ng	-0.01
79) Terphenyl-d14	13.09	244	1446420	101.46	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.90	88	142114	28.234	ng	90
3) Pyridine	3.68	79	416654	37.164	ng	# 85
4) n-Nitrosodimethylamine	3.63	42	214055	45.572	ng	90
6) Aniline	6.64	93	317813	19.854	ng	# 52
8) 2-Chlorophenol	6.76	128	440676	39.785	ng	99
9) Benzaldehyde	6.52	77	229847	29.989	ng	97
10) Phenol	6.61	94	691170	52.619	ng	# 68
11) bis(2-Chloroethyl)ether	6.70	93	409957	37.935	ng	93
12) 1,3-Dichlorobenzene	6.91	146	455880	37.400	ng	98
13) 1,4-Dichlorobenzene	6.99	146	468434	37.998	ng	99
14) 1,2-Dichlorobenzene	7.14	146	429656	37.543	ng	100
15) Benzyl Alcohol	7.11	79	398450	42.158	ng	92
16) 2,2'-oxybis(1-Chloropropan	7.23	45	676708	39.164	ng	68
17) 2-Methylphenol	7.22	107	370063	41.922	ng	# 80
18) Hexachloroethane	7.48	117	169816	37.624	ng	97
19) n-Nitroso-di-n-propylamine	7.38	70	303634	38.515	ng	97
20) 3+4-Methylphenols	7.37	107	467179	40.943	ng	91
22) Acetophenone	7.37	105	607606	41.897	ng	95
24) Nitrobenzene	7.55	77	466783	42.608	ng	97
25) Isophorone	7.79	82	850854	47.039	ng	96
26) 2-Nitrophenol	7.87	139	234113	44.907	ng	90
27) 2,4-Dimethylphenol	7.89	122	407072	47.096	ng	99
28) bis(2-Chloroethoxy)methane	7.99	93	534156	43.470	ng	99
29) 2,4-Dichlorophenol	8.10	162	379310	43.438	ng	99
30) 1,2,4-Trichlorobenzene	8.19	180	392045	40.081	ng	96
31) Naphthalene	8.27	128	1285194	44.305	ng	100
32) Benzoic acid	7.89	122	407072	60.936	ng	# 11
33) 4-Chloroaniline	8.32	127	135990	10.417	ng	99
34) Hexachlorobutadiene	8.38	225	208807	38.448	ng	98
35) Caprolactam	8.69	113	132922m	43.673	ng	
36) 4-Chloro-3-methylphenol	8.80	107	411808	43.611	ng	96
37) 2-Methylnaphthalene	8.97	142	877206	45.706	ng	98
38) 1-Methylnaphthalene	9.07	142	821483	44.292	ng	98
40) 1,2,4,5-Tetrachlorobenzene	9.13	216	361649	38.709	ng	98

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102518\
 Data File : BF110166.D
 Acq On : 25 Oct 2018 17:54
 Operator : JU/SJ
 Sample : J5636-02MS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HP-GRID-BMS

Manual Integrations
 APPROVED

Sohil
 10/26/2018 3:49:18 PM

Quant Time: Oct 26 04:22:55 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 23 15:06:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.12	237	308595	64.597	nq	98
43) 2,4,6-Trichlorophenol	9.25	196	257391	42.714	nq	96
44) 2,4,5-Trichlorophenol	9.28	196	272640	42.803	nq	94
46) 1,1'-Biphenyl	9.43	154	1041368	38.405	nq	99
47) 2-Chloronaphthalene	9.46	162	790265	39.732	nq	98
48) 2-Nitroaniline	9.55	65	267249	44.340	nq	97
49) Acenaphthylene	9.87	152	1254438	43.666	nq	98
50) Dimethylphthalate	9.73	163	1187123	53.633	nq	100
51) 2,6-Dinitrotoluene	9.80	165	212153	44.497	nq #	83
52) Acenaphthene	10.05	154	767854	40.403	nq	97
53) 3-Nitroaniline	9.96	138	143285	25.272	nq	93
54) 2,4-Dinitrophenol	10.07	184	21308	15.707	nq	88
55) Dibenzofuran	10.22	168	1081018	40.627	nq	96
56) 4-Nitrophenol	10.13	139	450390	107.330	nq	96
57) 2,4-Dinitrotoluene	10.20	165	276918	45.727	nq #	82
58) Fluorene	10.56	166	854098	43.129	nq	100
59) 2,3,4,6-Tetrachlorophenol	10.33	232	217440m	41.881	nq	
60) Diethylphthalate	10.43	149	906734	41.966	nq	99
61) 4-Chlorophenyl-phenylether	10.55	204	399153	38.208	nq	98
62) 4-Nitroaniline	10.59	138	214348	38.010	nq	91
63) Azobenzene	10.71	77	872069	40.662	nq	98
65) 4,6-Dinitro-2-methylphenol	10.61	198	60605	24.404	nq	93
66) n-Nitrosodiphenylamine	10.67	169	767465	43.045	nq	97
67) 4-Bromophenyl-phenylether	11.05	248	241777	41.006	nq	98
68) Hexachlorobenzene	11.11	284	243021	38.846	nq #	90
69) Atrazine	11.20	200	246736	43.791	nq	99
70) Pentachlorophenol	11.30	266	204455	56.046	nq	97
71) Phenanthrene	11.53	178	1356656	47.698	nq	99
72) Anthracene	11.58	178	1275685	44.747	nq	100
73) Carbazole	11.73	167	1064920	38.257	nq	100
74) Di-n-butylphthalate	12.05	149	1314607	41.816	nq	100
75) Fluoranthene	12.72	202	1284871	44.512	nq	99
77) Benzidine	12.84	184	290552	29.776	nq	96
78) Pyrene	12.96	202	1238938	58.287	nq	99
80) Butylbenzylphthalate	13.56	149	450153	48.792	nq	97
81) Benzo(a)anthracene	14.14	228	820044	46.640	nq	98
82) 3,3'-Dichlorobenzidine	14.10	252	148419	24.242	nq	98
83) Chrysene	14.18	228	758601	45.426	nq	99
84) Bis(2-ethylhexyl)phthalate	14.11	149	595301	47.733	nq	96
85) Di-n-octyl phthalate	14.73	149	886150	45.696	nq	99
86) Indeno(1,2,3-cd)pyrene	17.24	276	659439	47.599	nq	96
88) Benzo(b)fluoranthene	15.22	252	746342	46.975	nq	99
89) Benzo(k)fluoranthene	15.25	252	596814m	38.209	nq	
90) Benzo(a)pyrene	15.61	252	654761	44.786	nq	98
91) Dibenzo(a,h)anthracene	17.26	278	551954	43.755	nq	98
92) Benzo(g,h,i)perylene	17.72	276	515361	41.459	nq #	95

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

