

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF061218\
 Data File : BF106365.D
 Acq On : 13 Jun 2018 00:59
 Operator : JU/SJ
 Sample : PB110114BS
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB110114BS

Manual Integrations
 APPROVED

Sohil
 6/13/2018 4:56:29 PM

Quant Time: Jun 13 04:46:54 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 11 15:00:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.98	152	150703	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	531037	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	215809	20.00	ng	0.00
63) Phenanthrene-d10	11.51	188	398097	20.00	ng	0.00
75) Chrysene-d12	14.16	240	370101	20.00	ng	0.00
86) Perylene-d12	15.69	264	265003	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.62	112	1068555	120.01	ng	0.04
7) Phenol-d6	6.62	99	1317305	117.10	ng	0.04
23) Nitrobenzene-d5	7.54	82	838090	92.84	ng	0.00
41) 2,4,6-Tribromophenol	10.81	330	291962	140.61	ng	0.00
44) 2-Fluorobiphenyl	9.34	172	1229757	111.17	ng	0.00
78) Terphenyl-d14	13.09	244	1377022	92.41	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.89	88	150889	32.744	ng	96
3) Pyridine	3.65	79	448296	34.909	ng	99
4) n-Nitrosodimethylamine	3.62	42	208627	35.463	ng	98
6) Aniline	6.64	93	193787	11.797	ng	# 2
8) 2-Chlorophenol	6.78	128	371700	38.818	ng	99
9) Benzaldehyde	6.53	77	274290	32.145	ng	99
10) Phenol	6.64	94	480046	39.088	ng	78
11) bis(2-Chloroethyl)ether	6.71	93	418957	39.850	ng	98
12) 1,3-Dichlorobenzene	6.92	146	431290	38.270	ng	99
13) 1,4-Dichlorobenzene	7.00	146	435788	38.105	ng	99
14) 1,2-Dichlorobenzene	7.15	146	395906	36.774	ng	98
15) Benzyl Alcohol	7.12	79	333438	34.610	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.24	45	581513	38.586	ng	54
17) 2-Methylphenol	7.25	107	318629	37.297	ng	# 87
18) Hexachloroethane	7.49	117	133291	32.710	ng	91
19) n-Nitroso-di-n-propylamine	7.38	70	285824	33.866	ng	96
20) 3+4-Methylphenols	7.41	107	440682	40.336	ng	91
22) Acetophenone	7.38	105	531587	39.722	ng	# 98
24) Nitrobenzene	7.56	77	414150	42.212	ng	98
25) Isophorone	7.80	82	749739	42.523	ng	96
26) 2-Nitrophenol	7.87	139	187704	49.289	ng	95
27) 2,4-Dimethylphenol	7.92	122	333099	44.626	ng	98
28) bis(2-Chloroethoxy)methane	8.00	93	482345	42.190	ng	99
29) 2,4-Dichlorophenol	8.14	162	322637	44.804	ng	98
30) 1,2,4-Trichlorobenzene	8.20	180	324002	40.128	ng	95
31) Naphthalene	8.28	128	979055	40.784	ng	99
32) Benzoic acid	8.04	122	169915	33.466	ng	93
33) 4-Chloroaniline	8.33	127	90187	8.478	ng	96
34) Hexachlorobutadiene	8.39	225	219123	42.243	ng	98
35) Caprolactam	8.70	113	94530	38.810	ng	95
36) 4-Chloro-3-methylphenol	8.83	107	303419	38.091	ng	96
37) 2-Methylnaphthalene	8.97	142	650438	39.781	ng	96
39) 1,2,4,5-Tetrachlorobenzene	9.14	216	326829	46.651	ng	# 96
40) Hexachlorocyclopentadiene	9.12	237	41293	10.683	ng	93

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)

42) 2,4,6-Trichlorophenol	9.26	196	210863	47.230	nq	99
43) 2,4,5-Trichlorophenol	9.32	196	177371m	36.852	nq	
45) 1,1'-Biphenyl	9.44	154	741949	44.387	nq	99
46) 2-Chloronaphthalene	9.47	162	577366	43.162	nq	96
47) 2-Nitroaniline	9.56	65	198352	47.136	nq	89
48) Acenaphthylene	9.88	152	892143	43.573	nq	98
49) Dimethylphthalate	9.73	163	711229	46.129	nq	99
50) 2,6-Dinitrotoluene	9.80	165	140369	44.569	nq	90
51) Acenaphthene	10.05	154	527698	39.908	nq	99
52) 3-Nitroaniline	9.97	138	89815	24.027	nq	100
53) 2,4-Dinitrophenol	10.08	184	44878	36.308	nq #	17
54) Dibenzofuran	10.22	168	742709	41.712	nq	97
55) 4-Nitrophenol	10.18	139	200763	70.070	nq	91
56) 2,4-Dinitrotoluene	10.20	165	173548	43.881	nq	98
57) Fluorene	10.57	166	572271	40.565	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.35	232	167326	43.863	nq #	85
59) Diethylphthalate	10.43	149	655292	42.820	nq	95
60) 4-Chlorophenyl-phenylether	10.55	204	308324	41.553	nq	92
61) 4-Nitroaniline	10.58	138	128760	35.404	nq	98
62) Azobenzene	10.72	77	631410	39.568	nq	94
64) 4,6-Dinitro-2-methylphenol	10.61	198	40341	22.616	nq	90
65) n-Nitrosodiphenylamine	10.68	169	540667	40.410	nq	95
66) 4-Bromophenyl-phenylether	11.05	248	198659	43.875	nq #	85
67) Hexachlorobenzene	11.12	284	196208	42.032	nq #	84
68) Atrazine	11.20	200	177266	42.993	nq	96
69) Pentachlorophenol	11.32	266	174922	60.865	nq	97
70) Phenanthrene	11.54	178	825073	42.319	nq	98
71) Anthracene	11.59	178	855694	43.810	nq	99
72) Carbazole	11.74	167	774074	42.928	nq	98
73) Di-n-butylphthalate	12.06	149	929281	46.305	nq	99
74) Fluoranthene	12.72	202	956806	46.018	nq	98
76) Benzidine	12.85	184	686177	55.708	nq	99
77) Pyrene	12.96	202	978107	42.213	nq	97
79) Butylbenzylphthalate	13.57	149	445229	51.333	nq #	95
80) Benzo(a)anthracene	14.15	228	917112	41.641	nq	97
81) 3,3'-Dichlorobenzidine	14.11	252	317608	42.741	nq #	97
82) Chrysene	14.19	228	842387	41.892	nq	99
83) Bis(2-ethylhexyl)phthalate	14.12	149	606284	48.752	nq	100
84) Di-n-octyl phthalate	14.75	149	1001812	55.879	nq	99
85) Indeno(1,2,3-cd)pyrene	17.27	276	634916	39.563	nq	93
87) Benzo(b)fluoranthene	15.24	252	701635	43.800	nq	96
88) Benzo(k)fluoranthene	15.27	252	687037	43.306	nq #	97
89) Benzo(a)pyrene	15.63	252	637303	44.231	nq	97
90) Dibenzo(a,h)anthracene	17.29	278	537804	44.771	nq	96
91) Benzo(g,h,i)perylene	17.75	276	534232	45.763	nq #	93

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

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