

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100318\
 Data File : BF109578.D
 Acq On : 4 Oct 2018 6:20
 Operator : JU/SJ
 Sample : PB113461BSD
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB113461BSD

Manual Integrations
 APPROVED

Sohil
 10/4/2018 11:44:35 AM

Quant Time: Oct 04 07:29:55 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 22 13:32:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	88210	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	326588	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	141189	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	241594	20.00	ng	0.00
75) Chrysene-d12	13.84	240	192000	20.00	ng	0.00
86) Perylene-d12	15.24	264	159628	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.33	112	483407	90.26	ng	0.01
7) Phenol-d6	6.35	99	611492	89.48	ng	0.00
23) Nitrobenzene-d5	7.27	82	560026	101.23	ng	-0.01
41) 2,4,6-Tribromophenol	10.52	330	135963	97.69	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	1009479	107.83	ng	-0.01
78) Terphenyl-d14	12.79	244	857946	109.66	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.45	88	87115	35.319	ng	95
3) Pyridine	3.15	79	228747	36.033	ng	89
4) n-Nitrosodimethylamine	3.09	42	104670	38.744	ng	# 98
6) Aniline	6.37	93	251595	28.776	ng	# 31
8) 2-Chlorophenol	6.49	128	265302	44.547	ng	97
9) Benzaldehyde	6.25	77	113033	25.747	ng	99
10) Phenol	6.36	94	312750	40.411	ng	# 70
11) bis(2-Chloroethyl)ether	6.45	93	249783	41.247	ng	93
12) 1,3-Dichlorobenzene	6.65	146	286617	41.799	ng	97
13) 1,4-Dichlorobenzene	6.72	146	302383	43.772	ng	98
14) 1,2-Dichlorobenzene	6.87	146	266497	41.819	ng	99
15) Benzyl Alcohol	6.85	79	214136	40.936	ng	95
16) 2,2'-oxybis(1-Chloropropan	6.99	45	343499	39.989	ng	80
17) 2-Methylphenol	6.97	107	209806	43.539	ng	# 93
18) Hexachloroethane	7.22	117	89951	36.959	ng	96
19) n-Nitroso-di-n-propylamine	7.12	70	183185	40.908	ng	99
20) 3+4-Methylphenols	7.12	107	260454	44.767	ng	# 69
22) Acetophenone	7.12	105	372420	49.323	ng	98
24) Nitrobenzene	7.29	77	271764	46.199	ng	97
25) Isophorone	7.53	82	500561	50.245	ng	97
26) 2-Nitrophenol	7.60	139	116167	49.936	ng	99
27) 2,4-Dimethylphenol	7.65	122	228454	52.628	ng	99
28) bis(2-Chloroethoxy)methane	7.74	93	319636	48.468	ng	99
29) 2,4-Dichlorophenol	7.85	162	219990	48.235	ng	97
30) 1,2,4-Trichlorobenzene	7.93	180	229212	44.976	ng	98
31) Naphthalene	8.01	128	725628	47.547	ng	99
32) Benzoic acid	7.78	122	104414	37.181	ng	98
33) 4-Chloroaniline	8.06	127	162858	24.427	ng	99
34) Hexachlorobutadiene	8.13	225	139516	45.411	ng	97
35) Caprolactam	8.43	113	66431m	45.622	ng	
36) 4-Chloro-3-methylphenol	8.55	107	224071	46.688	ng	99
37) 2-Methylnaphthalene	8.70	142	483835	49.179	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.87	216	227291	50.584	ng	97
40) Hexachlorocyclopentadiene	8.85	237	50474	25.754	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.98	196	137738	47.761	nq	100
43) 2,4,5-Trichlorophenol	9.03	196	156650	51.432	nq	95
45) 1,1'-Biphenyl	9.16	154	579753	50.398	nq	98
46) 2-Chloronaphthalene	9.19	162	431839	46.990	nq	99
47) 2-Nitroaniline	9.28	65	139109	53.390	nq	93
48) Acenaphthylene	9.60	152	685496	49.445	nq	99
49) Dimethylphthalate	9.46	163	533785	51.979	nq	100
50) 2,6-Dinitrotoluene	9.52	165	104274	51.270	nq	90
51) Acenaphthene	9.77	154	374540	44.684	nq	99
52) 3-Nitroaniline	9.69	138	97611	41.443	nq	96
53) 2,4-Dinitrophenol	9.80	184	18033	30.848	nq #	38
54) Dibenzofuran	9.94	168	561644	45.056	nq	96
55) 4-Nitrophenol	9.87	139	132532	74.695	nq	94
56) 2,4-Dinitrotoluene	9.93	165	123150	47.855	nq #	93
57) Fluorene	10.28	166	428733	48.204	nq	97
58) 2,3,4,6-Tetrachlorophenol	10.06	232	123175	51.076	nq #	85
59) Diethylphthalate	10.16	149	510199	49.062	nq	98
60) 4-Chlorophenyl-phenylether	10.27	204	212849	45.818	nq	94
61) 4-Nitroaniline	10.30	138	108684	47.316	nq	95
62) Azobenzene	10.43	77	461695	42.733	nq	96
64) 4,6-Dinitro-2-methylphenol	10.34	198	17365	22.039	nq	85
65) n-Nitrosodiphenylamine	10.39	169	409706	51.328	nq	95
66) 4-Bromophenyl-phenylether	10.76	248	133422	49.732	nq #	87
67) Hexachlorobenzene	10.83	284	135336	47.498	nq #	88
68) Atrazine	10.92	200	132738	54.070	nq	98
69) Pentachlorophenol	11.03	266	129314	75.829	nq	99
70) Phenanthrene	11.24	178	589700	48.886	nq	99
71) Anthracene	11.29	178	631961	51.653	nq	99
72) Carbazole	11.44	167	551727	45.324	nq	100
73) Di-n-butylphthalate	11.78	149	708160	50.533	nq	99
74) Fluoranthene	12.42	202	626264	48.581	nq	96
76) Benzidine	12.54	184	482611	69.716	nq	98
77) Pyrene	12.64	202	620087	45.063	nq	100
79) Butylbenzylphthalate	13.27	149	282776	48.303	nq	95
80) Benzo(a)anthracene	13.83	228	523826	45.295	nq	99
81) 3,3'-Dichlorobenzidine	13.79	252	212540	48.358	nq	99
82) Chrysene	13.87	228	535048	46.678	nq	99
83) Bis(2-ethylhexyl)phthalate	13.83	149	357971	48.485	nq	97
84) Di-n-octyl phthalate	14.43	149	666039	52.761	nq	97
85) Indeno(1,2,3-cd)pyrene	16.59	276	396158	42.639	nq #	92
87) Benzo(b)fluoranthene	14.84	252	431565	49.201	nq	98
88) Benzo(k)fluoranthene	14.87	252	463898m	55.734	nq	
89) Benzo(a)pyrene	15.19	252	408701	51.709	nq	98
90) Dibenzo(a,h)anthracene	16.60	278	334622	51.605	nq #	96
91) Benzo(g,h,i)perylene	17.00	276	323585	50.776	nq #	91

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

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