

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122818\
 Data File : BF111800.D
 Acq On : 28 Dec 2018 20:29
 Operator : JU/SJ
 Sample : PB116072BSD
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB116072BSD

Manual Integrations
 APPROVED

Sohil
 1/2/2019 2:51:45 PM

Quant Time: Dec 29 02:52:33 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 19 17:39:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	139747	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	504027	20.00	ng	0.00
39) Acenaphthene-d10	9.95	164	235941	20.00	ng	0.01
64) Phenanthrene-d10	11.43	188	363024	20.00	ng	0.01
76) Chrysene-d12	14.07	240	305802	20.00	ng	0.02
87) Perylene-d12	15.56	264	242349	20.00	ng	0.04

System Monitoring Compounds

5) 2-Fluorophenol	5.55	112	998320	121.77	ng	0.04
7) Phenol-d6	6.55	99	1274493	122.15	ng	0.03
23) Nitrobenzene-d5	7.46	82	814188	94.03	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	323198	115.93	ng	0.01
45) 2-Fluorobiphenyl	9.26	172	1532004	94.64	ng	0.00
79) Terphenyl-d14	13.01	244	1277432	79.22	ng	0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.78	88	127694	33.104	ng	95
3) Pyridine	3.53	79	372130	37.030	ng	97
4) n-Nitrosodimethylamine	3.48	42	213665	43.444	ng	# 79
6) Aniline	6.57	93	274318	20.625	ng	# 6
8) 2-Chlorophenol	6.70	128	383288	42.203	ng	99
9) Benzaldehyde	6.45	77	201376	28.062	ng	97
10) Phenol	6.56	94	452901	38.471	ng	77
11) bis(2-Chloroethyl)ether	6.64	93	358467	40.634	ng	99
12) 1,3-Dichlorobenzene	6.85	146	443143	42.104	ng	98
13) 1,4-Dichlorobenzene	6.92	146	442469	41.452	ng	98
14) 1,2-Dichlorobenzene	7.08	146	415143	41.686	ng	96
15) Benzyl Alcohol	7.05	79	346741	41.844	ng	93
16) 2,2'-oxybis(1-Chloropropan	7.17	45	587511	36.160	ng	69
17) 2-Methylphenol	7.17	107	306556	42.155	ng	# 88
18) Hexachloroethane	7.42	117	143420	39.873	ng	# 83
19) n-Nitroso-di-n-propylamine	7.32	70	273096	39.412	ng	99
20) 3+4-Methylphenols	7.32	107	395129	43.001	ng	# 82
22) Acetophenone	7.31	105	523870	41.027	ng	# 91
24) Nitrobenzene	7.49	77	408287	44.988	ng	98
25) Isophorone	7.72	82	719057	46.776	ng	97
26) 2-Nitrophenol	7.80	139	190143	51.659	ng	# 87
27) 2,4-Dimethylphenol	7.85	122	335776	46.277	ng	94
28) bis(2-Chloroethoxy)methane	7.93	93	448568	43.850	ng	98
29) 2,4-Dichlorophenol	8.05	162	336995	43.181	ng	96
30) 1,2,4-Trichlorobenzene	8.13	180	360738	41.480	ng	97
31) Naphthalene	8.21	128	1100833	45.456	ng	96
32) Benzoic acid	7.95	122	170814	35.606	ng	96
33) 4-Chloroaniline	8.25	127	112175	10.717	ng	99
34) Hexachlorobutadiene	8.33	225	221512	41.792	ng	98
35) Caprolactam	8.62	113	99510m	37.629	ng	
36) 4-Chloro-3-methylphenol	8.75	107	326263	42.529	ng	100
37) 2-Methylnaphthalene	8.90	142	736953	46.772	ng	100
38) 1-Methylnaphthalene	9.00	142	690985	45.662	ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.07	216	353494	45.595	ng	# 97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	9.06	237	193590	51.456	nq	100
43) 2,4,6-Trichlorophenol	9.18	196	230175	44.467	nq	98
44) 2,4,5-Trichlorophenol	9.23	196	235204	44.292	nq	91
46) 1,1'-Biphenyl	9.36	154	893221	44.847	nq	95
47) 2-Chloronaphthalene	9.39	162	678620	43.005	nq	93
48) 2-Nitroaniline	9.48	65	207948	50.655	nq	99
49) Acenaphthylene	9.80	152	1037890	45.048	nq	97
50) Dimethylphthalate	9.66	163	819720	47.510	nq	98
51) 2,6-Dinitrotoluene	9.72	165	171527	49.857	nq	98
52) Acenaphthene	9.98	154	624989	43.982	nq	97
53) 3-Nitroaniline	9.89	138	98690	26.897	nq	98
54) 2,4-Dinitrophenol	9.99	184	56032	53.421	nq	97
55) Dibenzofuran	10.15	168	913633	42.557	nq	94
56) 4-Nitrophenol	10.07	139	234087	83.599	nq	88
57) 2,4-Dinitrotoluene	10.12	165	208342	50.387	nq	95
58) Fluorene	10.49	166	677002	43.762	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.27	232	178342	39.907	nq	90
60) Diethylphthalate	10.36	149	773597	45.708	nq	99
61) 4-Chlorophenyl-phenylether	10.48	204	357517	41.830	nq	95
62) 4-Nitroaniline	10.50	138	137185	38.469	nq	89
63) Azobenzene	10.64	77	668516	39.345	nq	96
65) 4,6-Dinitro-2-methylphenol	10.53	198	41912	30.733	nq	81
66) n-Nitrosodiphenylamine	10.60	169	598871	49.144	nq	98
67) 4-Bromophenyl-phenylether	10.97	248	214965	47.742	nq	97
68) Hexachlorobenzene	11.05	284	224355	44.689	nq	95
69) Atrazine	11.12	200	193873	45.795	nq	96
70) Pentachlorophenol	11.24	266	199059	64.137	nq	99
71) Phenanthrene	11.46	178	892913	46.379	nq	94
72) Anthracene	11.50	178	918009	47.505	nq	95
73) Carbazole	11.66	167	773090	41.206	nq	95
74) Di-n-butylphthalate	11.99	149	1013518	48.181	nq	96
75) Fluoranthene	12.65	202	865518	44.395	nq	94
77) Benzidine	12.76	184	527638	45.853	nq	97
78) Pyrene	12.87	202	874670	40.503	nq	96
80) Butylbenzylphthalate	13.48	149	387080	43.973	nq	92
81) Benzo(a)anthracene	14.06	228	835643	47.882	nq	97
82) 3,3'-Dichlorobenzidine	14.02	252	294056	42.645	nq	# 98
83) Chrysene	14.10	228	774729	45.167	nq	# 96
84) Bis(2-ethylhexyl)phthalate	14.05	149	532959	48.066	nq	# 99
85) Di-n-octyl phthalate	14.66	149	908118	52.862	nq	# 95
86) Indeno(1,2,3-cd)pyrene	17.07	276	574981	39.432	nq	# 88
88) Benzo(b)fluoranthene	15.12	252	732097	51.688	nq	# 95
89) Benzo(k)fluoranthene	15.15	252	640588	47.506	nq	# 97
90) Benzo(a)pyrene	15.50	252	621903	48.330	nq	# 94
91) Dibenzo(a,h)anthracene	17.08	278	489295	43.423	nq	# 90
92) Benzo(g,h,i)perylene	17.52	276	461275	41.923	nq	# 85

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

