

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090418\
 Data File : BF108761.D
 Acq On : 5 Sep 2018 2:52
 Operator : JU/SJ
 Sample : PB112567BS
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB112567BS

Manual Integrations
 APPROVED

Sohil
 9/5/2018 12:48:03 PM

Quant Time: Sep 05 08:21:22 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 30 15:06:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.00	152	68651	20.00	ng	-0.01
21) Naphthalene-d8	8.27	136	246641	20.00	ng	-0.01
38) Acenaphthene-d10	10.04	164	114976	20.00	ng	-0.01
63) Phenanthrene-d10	11.53	188	235491	20.00	ng	-0.02
75) Chrysene-d12	14.19	240	177342	20.00	ng	-0.01
86) Perylene-d12	15.74	264	126219	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.63	112	430175	108.86	ng	-0.02
7) Phenol-d6	6.62	99	577103	101.96	ng	-0.02
23) Nitrobenzene-d5	7.56	82	403389	83.02	ng	-0.01
41) 2,4,6-Tribromophenol	10.83	330	161837	106.62	ng	-0.02
44) 2-Fluorobiphenyl	9.34	172	762246	89.14	ng	-0.02
78) Terphenyl-d14	13.11	244	894243	97.97	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.92	88	64429	35.081	ng	95
3) Pyridine	3.72	79	132293	31.177	ng	# 82
4) n-Nitrosodimethylamine	3.70	42	48541	22.933	ng	85
6) Aniline	6.68	93	140508	23.006	ng	# 80
8) 2-Chlorophenol	6.78	128	190029	39.328	ng	95
9) Benzaldehyde	6.56	77	107438m	30.711	ng	
10) Phenol	6.64	94	187142	28.354	ng	81
11) bis(2-Chloroethyl)ether	6.72	93	212608	36.581	ng	97
12) 1,3-Dichlorobenzene	6.93	146	201937	36.045	ng	98
13) 1,4-Dichlorobenzene	7.01	146	238865	39.060	ng	99
14) 1,2-Dichlorobenzene	7.16	146	230907	41.113	ng	97
15) Benzyl Alcohol	7.14	79	106971m	26.273	ng	
16) 2,2'-oxybis(1-Chloropropan	7.25	45	351764	39.453	ng	76
17) 2-Methylphenol	7.24	107	129439	34.455	ng	# 72
18) Hexachloroethane	7.50	117	67302	32.001	ng	93
19) n-Nitroso-di-n-propylamine	7.39	70	141171	37.245	ng	91
20) 3+4-Methylphenols	7.40	107	144065	30.607	ng	# 18
22) Acetophenone	7.40	105	275655	44.770	ng	# 88
24) Nitrobenzene	7.58	77	183308	37.205	ng	94
25) Isophorone	7.80	82	371286	45.998	ng	95
26) 2-Nitrophenol	7.90	139	63230	28.349	ng	# 88
27) 2,4-Dimethylphenol	7.92	122	161909	49.913	ng	96
28) bis(2-Chloroethoxy)methane	8.01	93	210752	41.332	ng	99
29) 2,4-Dichlorophenol	8.14	162	145782	37.882	ng	99
30) 1,2,4-Trichlorobenzene	8.21	180	173615	40.344	ng	95
31) Naphthalene	8.30	128	556115	47.338	ng	99
32) Benzoic acid	8.05	122	28784m	13.837	ng	
33) 4-Chloroaniline	8.36	127	139263	26.854	ng	93
34) Hexachlorobutadiene	8.40	225	119730	42.373	ng	99
35) Caprolactam	8.71	113	37512	36.581	ng	94
36) 4-Chloro-3-methylphenol	8.83	107	150251	36.377	ng	100
37) 2-Methylnaphthalene	8.99	142	342375	43.075	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.15	216	181577	44.484	ng	# 96
40) Hexachlorocyclopentadiene	9.12	237	12547	13.684	ng	96

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.27	196	89294	36.941	ng	99
43) 2,4,5-Trichlorophenol	9.32	196	129071	45.791	ng	95
45) 1,1'-Biphenyl	9.45	154	435085	43.468	ng	99
46) 2-Chloronaphthalene	9.48	162	333592	42.571	ng	98
47) 2-Nitroaniline	9.59	65	111360	42.490	ng	86
48) Acenaphthylene	9.90	152	514051	44.808	ng	99
49) Dimethylphthalate	9.74	163	393979	43.515	ng	# 99
50) 2,6-Dinitrotoluene	9.82	165	73927	37.043	ng	# 90
51) Acenaphthene	10.07	154	300366	43.789	ng	100
52) 3-Nitroaniline	10.01	138	95671	44.314	ng	91
54) Dibenzofuran	10.24	168	445581	41.514	ng	96
55) 4-Nitrophenol	10.19	139	48100	37.605	ng	# 81
56) 2,4-Dinitrotoluene	10.23	165	89128	33.729	ng	# 89
57) Fluorene	10.59	166	366075	44.010	ng	96
58) 2,3,4,6-Tetrachlorophenol	10.36	232	98220	38.897	ng	# 83
59) Diethylphthalate	10.44	149	384481	42.150	ng	96
60) 4-Chlorophenyl-phenylether	10.57	204	189507	40.149	ng	91
61) 4-Nitroaniline	10.63	138	89694	42.934	ng	93
62) Azobenzene	10.73	77	370933	40.912	ng	94
64) 4,6-Dinitro-2-methylphenol	10.65	198	2367	2.265	ng	# 1
65) n-Nitrosodiphenylamine	10.70	169	324204	41.438	ng	96
66) 4-Bromophenyl-phenylether	11.07	248	122268	42.245	ng	# 85
67) Hexachlorobenzene	11.14	284	126671	40.624	ng	# 84
68) Atrazine	11.21	200	110234	57.484	ng	96
69) Pentachlorophenol	11.34	266	87438	52.772	ng	99
70) Phenanthrene	11.55	178	516688	43.469	ng	99
71) Anthracene	11.61	178	621336	49.749	ng	99
72) Carbazole	11.77	167	518744	42.955	ng	99
73) Di-n-butylphthalate	12.07	149	614485	44.226	ng	99
74) Fluoranthene	12.75	202	637128	47.619	ng	93
76) Benzidine	12.88	184	492472	98.653	ng	99
77) Pyrene	12.98	202	610787	45.816	ng	100
79) Butylbenzylphthalate	13.58	149	268595	47.915	ng	97
80) Benzo(a)anthracene	14.18	228	457003	42.271	ng	99
81) 3,3'-Dichlorobenzidine	14.14	252	186421	47.547	ng	# 96
82) Chrysene	14.21	228	484226	46.574	ng	98
83) Bis(2-ethylhexyl)phthalate	14.12	149	361800	50.213	ng	100
84) Di-n-octyl phthalate	14.75	149	545702	51.511	ng	97
85) Indeno(1,2,3-cd)pyrene	17.36	276	323075	44.690	ng	# 89
87) Benzo(b)fluoranthene	15.27	252	324370	41.835	ng	# 97
88) Benzo(k)fluoranthene	15.30	252	364374	47.440	ng	# 96
89) Benzo(a)pyrene	15.67	252	320666	45.764	ng	# 97
90) Dibenzo(a,h)anthracene	17.37	278	279173	49.886	ng	# 92
91) Benzo(g,h,i)perylene	17.85	276	265306	48.778	ng	# 88

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

