

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102219\
 Data File : BF117475.D
 Acq On : 22 Oct 2019 13:46
 Operator : JU
 Sample : PB124054BSD
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB124054BSD

Manual Integrations
 APPROVED

mohammad
 10/23/2019 8:05:59 AM

Quant Time: Oct 22 14:57:38 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 18:53:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	43650	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	197543	20.00	ng	0.00
39) Acenaphthene-d10	9.77	164	105285	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	167490	20.00	ng	0.00
76) Chrysene-d12	13.90	240	118709	20.00	ng	0.00
87) Perylene-d12	15.33	264	97027	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	361440	123.19	ng	0.02
7) Phenol-d6	6.39	99	474388	120.38	ng	0.00
23) Nitrobenzene-d5	7.30	82	250328	62.56	ng	0.00
42) 2,4,6-Tribromophenol	10.56	330	101266	111.16	ng	0.00
45) 2-Fluorobiphenyl	9.09	172	456612	61.11	ng	0.00
79) Terphenyl-d14	12.83	244	446679	61.38	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.52	88	51304	41.589	ng	# 96
3) Pyridine	3.26	79	115301	38.181	ng	97
4) n-Nitrosodimethylamine	3.19	42	50666	46.247	ng	98
6) Aniline	6.40	93	161117	34.825	ng	# 27
8) 2-Chlorophenol	6.53	128	146215	47.226	ng	95
9) Benzaldehyde	6.29	77	93234	52.543	ng	97
10) Phenol	6.40	94	185823	46.201	ng	91
11) bis(2-Chloroethyl)ether	6.47	93	135979	45.707	ng	99
12) 1,3-Dichlorobenzene	6.67	146	148774	43.061	ng	98
13) 1,4-Dichlorobenzene	6.75	146	150273	42.875	ng	94
14) 1,2-Dichlorobenzene	6.90	146	144285	43.672	ng	98
15) Benzyl Alcohol	6.89	79	132078	46.911	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.01	45	138525	43.164	ng	70
17) 2-Methylphenol	7.00	107	122465	48.084	ng	96
18) Hexachloroethane	7.24	117	57807	42.256	ng	93
19) n-Nitroso-di-n-propylamine	7.15	70	110580	45.947	ng	96
20) 3+4-Methylphenols	7.16	107	158508	47.468	ng	# 81
22) Acetophenone	7.15	105	214323	40.858	ng	96
24) Nitrobenzene	7.32	77	167942	43.820	ng	95
25) Isophorone	7.56	82	306560	44.287	ng	99
26) 2-Nitrophenol	7.64	139	76412	44.703	ng	97
27) 2,4-Dimethylphenol	7.68	122	128063	50.131	ng	98
28) bis(2-Chloroethoxy)methane	7.77	93	181234	42.430	ng	99
29) 2,4-Dichlorophenol	7.89	162	120228	44.871	ng	99
30) 1,2,4-Trichlorobenzene	7.96	180	119615	41.600	ng	99
31) Naphthalene	8.04	128	423126	43.213	ng	99
32) Benzoic acid	7.84	122	69459	39.697	ng	95
33) 4-Chloroaniline	8.09	127	108015	26.139	ng	98
34) Hexachlorobutadiene	8.15	225	67909	40.804	ng	98
35) Caprolactam	8.47	113	37464m	43.525	ng	
36) 4-Chloro-3-methylphenol	8.59	107	138891	45.451	ng	100
37) 2-Methylnaphthalene	8.73	142	290200	43.984	ng	98
38) 1-Methylnaphthalene	8.83	142	275310	44.406	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.90	216	111017	39.149	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.88	237	63787	81.814	nq	98
43) 2,4,6-Trichlorophenol	9.02	196	74849	42.149	nq	98
44) 2,4,5-Trichlorophenol	9.06	196	79151	43.986	nq	96
46) 1,1'-Biphenyl	9.19	154	342876	39.338	nq	98
47) 2-Chloronaphthalene	9.22	162	263399	40.080	nq	100
48) 2-Nitroaniline	9.32	65	87628	40.617	nq	92
49) Acenaphthylene	9.63	152	413743	42.862	nq	99
50) Dimethylphthalate	9.49	163	313931	41.826	nq	100
51) 2,6-Dinitrotoluene	9.56	165	67910	43.358	nq	95
52) Acenaphthene	9.80	154	245667	41.481	nq	98
53) 3-Nitroaniline	9.73	138	50520	26.584	nq	92
54) 2,4-Dinitrophenol	9.86	184	45233	70.383	nq	93
55) Dibenzofuran	9.97	168	365578	42.606	nq	98
56) 4-Nitrophenol	9.92	139	71425	81.565	nq	86
57) 2,4-Dinitrotoluene	9.97	165	89568	43.012	nq	94
58) Fluorene	10.32	166	299440	42.201	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.10	232	63123	44.632	nq	99
60) Diethylphthalate	10.19	149	323581	42.623	nq	98
61) 4-Chlorophenyl-phenylether	10.30	204	131514	41.312	nq	99
62) 4-Nitroaniline	10.36	138	64495	36.006	nq	99
63) Azobenzene	10.47	77	339404	43.672	nq	99
65) 4,6-Dinitro-2-methylphenol	10.39	198	37232	43.722	nq	94
66) n-Nitrosodiphenylamine	10.43	169	261851	42.765	nq	97
67) 4-Bromophenyl-phenylether	10.79	248	75135	41.648	nq	99
68) Hexachlorobenzene	10.87	284	77555	40.141	nq	# 86
69) Atrazine	10.96	200	83082	53.331	nq	98
70) Pentachlorophenol	11.07	266	62670	108.202	nq	98
71) Phenanthrene	11.28	178	409816	42.138	nq	98
72) Anthracene	11.33	178	432867	43.378	nq	99
73) Carbazole	11.49	167	388871	42.642	nq	99
74) Di-n-butylphthalate	11.81	149	528846	43.401	nq	100
75) Fluoranthene	12.47	202	417289	42.860	nq	98
77) Benzidine	12.59	184	190869	74.739	nq	99
78) Pyrene	12.70	202	420556	41.245	nq	99
80) Butylbenzylphthalate	13.30	149	213681	42.113	nq	98
81) Benzo(a)anthracene	13.89	228	334149	41.730	nq	99
82) 3,3'-Dichlorobenzidine	13.84	252	81839	32.571	nq	98
83) Chrysene	13.93	228	327723	41.760	nq	98
84) Bis(2-ethylhexyl)phthalate	13.87	149	315736	44.422	nq	96
85) Di-n-octyl phthalate	14.48	149	511974	46.849	nq	100
86) Indeno(1,2,3-cd)pyrene	16.74	276	227522	38.832	nq	99
88) Benzo(b)fluoranthene	14.92	252	254724	44.458	nq	98
89) Benzo(k)fluoranthene	14.95	252	271775	46.280	nq	97
90) Benzo(a)pyrene	15.27	252	226474	43.475	nq	99
91) Dibenzo(a,h)anthracene	16.74	278	194733	43.232	nq	98
92) Benzo(g,h,i)perylene	17.16	276	184529	43.297	nq	97

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

