

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032620\
 Data File : BF119549.D
 Acq On : 27 Mar 2020 00:15
 Operator : CG/JU
 Sample : L2028-01MSD
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-30-(8-8.5)MSD

Manual Integrations
 APPROVED

mohammad
 3/30/2020 8:50:57 AM

Quant Time: Mar 27 05:04:11 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 26 16:30:13 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	286288	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	1136093	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	602853	20.00	ng	0.00
64) Phenanthrene-d10	11.37	188	1166985	20.00	ng	0.00
76) Chrysene-d12	14.02	240	810479	20.00	ng	0.00
87) Perylene-d12	15.48	264	771364	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	1499870	83.40	ng	0.01
7) Phenol-d6	6.46	99	1993122	86.76	ng	0.00
23) Nitrobenzene-d5	7.40	82	1231405	58.91	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	612104	98.42	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	2401750	59.02	ng	0.00
79) Terphenyl-d14	12.96	244	2811425	67.96	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.63	88	315454	35.516	ng	99
3) Pyridine	3.36	79	849500	34.716	ng	96
4) n-Nitrosodimethylamine	3.32	42	501429	42.021	ng	95
6) Aniline	6.50	93	1137419	35.873	ng	100
8) 2-Chlorophenol	6.62	128	988484	52.333	ng	98
9) Benzaldehyde	6.38	77	608073	41.724	ng	98
10) Phenol	6.47	94	1279238	52.186	ng	98
11) bis(2-Chloroethyl)ether	6.57	93	947196	48.313	ng	99
12) 1,3-Dichlorobenzene	6.77	146	963123	47.113	ng	99
13) 1,4-Dichlorobenzene	6.85	146	986214	48.230	ng	99
14) 1,2-Dichlorobenzene	7.01	146	926189	48.459	ng	98
15) Benzyl Alcohol	6.98	79	835701	48.360	ng	95
16) 2,2'-oxybis(1-Chloropropan	7.12	45	1799597	45.507	ng	97
17) 2-Methylphenol	7.09	107	802364	46.363	ng	98
18) Hexachloroethane	7.35	117	373744	49.876	ng	98
19) n-Nitroso-di-n-propylamine	7.26	70	691505	49.939	ng	96
20) 3+4-Methylphenols	7.25	107	965502	45.523	ng	# 79
22) Acetophenone	7.25	105	1272692	46.882	ng	99
24) Nitrobenzene	7.42	77	1035375	46.346	ng	95
25) Isophorone	7.66	82	1963213	46.297	ng	99
26) 2-Nitrophenol	7.73	139	536076	60.945	ng	97
27) 2,4-Dimethylphenol	7.77	122	836288	45.980	ng	99
28) bis(2-Chloroethoxy)methane	7.87	93	1232000	49.132	ng	99
29) 2,4-Dichlorophenol	7.98	162	832553	49.818	ng	97
30) 1,2,4-Trichlorobenzene	8.06	180	873906	47.284	ng	99
31) Naphthalene	8.15	128	2720596	46.534	ng	99
32) Benzoic acid	7.91	122	679345	58.065	ng	97
33) 4-Chloroaniline	8.19	127	413085	15.420	ng	99
34) Hexachlorobutadiene	8.26	225	498905	48.247	ng	98
35) Caprolactam	8.57	113	300465m	54.935	ng	
36) 4-Chloro-3-methylphenol	8.67	107	911049	49.878	ng	98
37) 2-Methylnaphthalene	8.84	142	1883602	49.091	ng	98
38) 1-Methylnaphthalene	8.94	142	1797210	49.486	ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.00	216	809015	45.784	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.99	237	998428	99.389	nq	97
43) 2,4,6-Trichlorophenol	9.12	196	638273	50.476	nq	98
44) 2,4,5-Trichlorophenol	9.16	196	648914	45.810	nq	97
46) 1,1'-Biphenyl	9.30	154	2247980	45.784	nq	99
47) 2-Chloronaphthalene	9.33	162	1791954	46.593	nq	99
48) 2-Nitroaniline	9.42	65	662076	49.874	nq	98
49) Acenaphthylene	9.75	152	2768765	45.843	nq	99
50) Dimethylphthalate	9.61	163	2419152	56.495	nq	99
51) 2,6-Dinitrotoluene	9.67	165	483776	52.708	nq	94
52) Acenaphthene	9.92	154	1974672	52.445	nq	99
53) 3-Nitroaniline	9.83	138	272937	23.554	nq	99
54) 2,4-Dinitrophenol	9.95	184	514413	125.023	nq	91
55) Dibenzofuran	10.09	168	2514403	46.578	nq	98
56) 4-Nitrophenol	10.00	139	867832	94.937	nq	98
57) 2,4-Dinitrotoluene	10.07	165	643914	55.691	nq	# 89
58) Fluorene	10.43	166	1943478	48.684	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.21	232	554769	52.394	nq	98
60) Diethylphthalate	10.32	149	2209849	50.847	nq	98
61) 4-Chlorophenyl-phenylether	10.43	204	944785	48.397	nq	97
62) 4-Nitroaniline	10.46	138	525179	47.264	nq	99
63) Azobenzene	10.59	77	2093785	47.841	nq	98
65) 4,6-Dinitro-2-methylphenol	10.48	198	339165	69.310	nq	88
66) n-Nitrosodiphenylamine	10.55	169	1812108	47.301	nq	99
67) 4-Bromophenyl-phenylether	10.92	248	630728	47.457	nq	98
68) Hexachlorobenzene	10.98	284	674669	49.599	nq	99
69) Atrazine	11.08	200	570962	54.363	nq	96
70) Pentachlorophenol	11.17	266	767907	96.279	nq	100
71) Phenanthrene	11.40	178	2759694	45.606	nq	98
72) Anthracene	11.45	178	2867409	46.625	nq	98
73) Carbazole	11.60	167	2681223	44.046	nq	99
74) Di-n-butylphthalate	11.94	149	3392477	52.097	nq	99
75) Fluoranthene	12.59	202	3068097	46.850	nq	98
77) Benzidine	12.71	184	1782192	51.959	nq	97
78) Pyrene	12.82	202	3083641	46.360	nq	99
80) Butylbenzylphthalate	13.44	149	1501267	55.804	nq	99
81) Benzo(a)anthracene	14.01	228	2386935	45.289	nq	99
82) 3,3'-Dichlorobenzidine	13.97	252	548523	26.396	nq	97
83) Chrysene	14.04	228	2535406	46.613	nq	98
84) Bis(2-ethylhexyl)phthalate	14.00	149	1940085	60.495	nq	99
85) Di-n-octyl phthalate	14.62	149	3410138	60.462	nq	98
86) Indeno(1,2,3-cd)pyrene	16.95	276	2302458	44.946	nq	97
88) Benzo(b)fluoranthene	15.06	252	2506408	48.643	nq	99
89) Benzo(k)fluoranthene	15.09	252	2343043	47.755	nq	100
90) Benzo(a)pyrene	15.42	252	2153400	45.986	nq	99
91) Dibenzo(a,h)anthracene	16.97	278	1884006	45.998	nq	98
92) Benzo(g,h,i)perylene	17.40	276	1869699	45.439	nq	96

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

