

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF081420\
 Data File : BF121290.D
 Acq On : 14 Aug 2020 17:46
 Operator : JU/CG
 Sample : PB130964BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB130964BS

Manual Integrations
 APPROVED

mohammad
 8/18/2020 1:42:03 PM

Quant Time: Aug 14 18:19:04 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF081320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 13 15:23:26 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	67303	20.00	ng	-0.02
21) Naphthalene-d8	7.99	136	266433	20.00	ng	-0.01
39) Acenaphthene-d10	9.75	164	141954	20.00	ng	-0.01
64) Phenanthrene-d10	11.23	188	288312	20.00	ng	-0.01
76) Chrysene-d12	13.86	240	264486	20.00	ng	-0.01
86) Perylene-d12	15.27	264	310930	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	478887	107.87	ng	0.00
7) Phenol-d6	6.35	99	571750	99.94	ng	-0.01
23) Nitrobenzene-d5	7.27	82	363826	74.63	ng	-0.02
42) 2,4,6-Tribromophenol	10.53	330	187769	107.44	ng	-0.01
45) 2-Fluorobiphenyl	9.07	172	729434	71.08	ng	-0.01
79) Terphenyl-d14	12.81	244	875518	64.53	ng	-0.01

Target Compounds					Qvalue
2) 1,4-Dioxane	2.29	88	68606	36.315 ng	98
3) Pyridine	3.00	79	154676	31.452 ng	96
4) n-Nitrosodimethylamine	2.94	42	88809	44.094 ng	97
6) Aniline	6.36	93	253588	37.394 ng	99
8) 2-Chlorophenol	6.49	128	203108	41.607 ng	99
9) Benzaldehyde	6.25	77	124676	42.291 ng	99
10) Phenol	6.36	94	258351	41.631 ng	98
11) bis(2-Chloroethyl)ether	6.44	93	190916	42.620 ng	96
12) 1,3-Dichlorobenzene	6.65	146	207886	40.304 ng	99
13) 1,4-Dichlorobenzene	6.72	146	209769	40.237 ng	99
14) 1,2-Dichlorobenzene	6.87	146	201142	40.880 ng	100
15) Benzyl Alcohol	6.85	79	159729	41.682 ng	99
16) 2,2'-oxybis(1-Chloropropan	6.99	45	267079	41.406 ng	98
17) 2-Methylphenol	6.97	107	165984	43.046 ng	98
18) Hexachloroethane	7.22	117	76667	41.585 ng	98
19) n-Nitroso-di-n-propylamine	7.12	70	128281	40.359 ng	97
20) 3+4-Methylphenols	7.12	107	200840	41.053 ng	# 84
22) Acetophenone	7.12	105	266629	38.854 ng	# 98
24) Nitrobenzene	7.29	77	204940	38.743 ng	94
25) Isophorone	7.53	82	372804	38.089 ng	96
26) 2-Nitrophenol	7.61	139	106315	41.481 ng	98
27) 2,4-Dimethylphenol	7.65	122	178387	43.440 ng	98
28) bis(2-Chloroethoxy)methane	7.75	93	238835	43.626 ng	100
29) 2,4-Dichlorophenol	7.85	162	166631	39.680 ng	98
30) 1,2,4-Trichlorobenzene	7.93	180	178570	38.528 ng	99
31) Naphthalene	8.01	128	572465	38.191 ng	100
32) Benzoic acid	7.76	122	66312	32.167 ng	97
33) 4-Chloroaniline	8.06	127	209059	34.176 ng	100
34) Hexachlorobutadiene	8.13	225	103697	36.978 ng	98
35) Caprolactam	8.42	113	55834m	44.142 ng	
36) 4-Chloro-3-methylphenol	8.55	107	176101	41.103 ng	97
37) 2-Methylnaphthalene	8.70	142	383955	38.926 ng	99
38) 1-Methylnaphthalene	8.80	142	368384	39.444 ng	100
40) 1,2,4,5-Tetrachlorobenzene	8.87	216	182135	38.292 ng	99

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41) Hexachlorocyclopentadiene	8.86	237	159800	78.274	nq	100
43) 2,4,6-Trichlorophenol	8.99	196	119797	39.602	nq	98
44) 2,4,5-Trichlorophenol	9.03	196	131244	39.945	nq	96
46) 1,1'-Biphenyl	9.17	154	476065	38.813	nq	98
47) 2-Chloronaphthalene	9.19	162	364313	38.585	nq	100
48) 2-Nitroaniline	9.29	65	110657	41.695	nq	98
49) Acenaphthylene	9.60	152	590280	39.361	nq	99
50) Dimethylphthalate	9.47	163	435507	40.771	nq	99
51) 2,6-Dinitrotoluene	9.53	165	96946	40.714	nq	94
52) Acenaphthene	9.77	154	340313	38.328	nq	100
53) 3-Nitroaniline	9.70	138	88688	32.621	nq	99
54) 2,4-Dinitrophenol	9.81	184	79658	65.994	nq	# 81
55) Dibenzofuran	9.95	168	533327	37.719	nq	98
56) 4-Nitrophenol	9.87	139	163354	93.104	nq	93
57) 2,4-Dinitrotoluene	9.94	165	131248	42.001	nq	98
58) Fluorene	10.29	166	406474	37.977	nq	100
59) 2,3,4,6-Tetrachlorophenol	10.07	232	112250	43.209	nq	99
60) Diethylphthalate	10.17	149	444494	40.921	nq	99
61) 4-Chlorophenyl-phenylether	10.29	204	199948	38.410	nq	98
62) 4-Nitroaniline	10.31	138	117154	42.979	nq	97
63) Azobenzene	10.45	77	423973	39.064	nq	99
65) 4,6-Dinitro-2-methylphenol	10.35	198	66442	38.533	nq	98
66) n-Nitrosodiphenylamine	10.40	169	381429	38.583	nq	100
67) 4-Bromophenyl-phenylether	10.77	248	130056	38.672	nq	97
68) Hexachlorobenzene	10.84	284	147689	38.511	nq	98
69) Atrazine	10.93	200	110407	37.338	nq	97
70) Pentachlorophenol	11.04	266	152357	88.498	nq	99
71) Phenanthrene	11.25	178	635496	38.529	nq	100
72) Anthracene	11.30	178	660786	39.922	nq	100
73) Carbazole	11.46	167	628080	38.800	nq	99
74) Di-n-butylphthalate	11.79	149	756044	41.378	nq	100
75) Fluoranthene	12.44	202	753043	41.002	nq	99
77) Benzidine	12.56	184	494574	71.131	nq	99
78) Pyrene	12.66	202	776793	39.440	nq	99
80) Butylbenzylphthalate	13.29	149	347315	43.171	nq	99
81) Benzo(a)anthracene	13.85	228	693772	38.674	nq	100
82) 3,3'-Dichlorobenzidine	13.82	252	220219	30.941	nq	99
83) Chrysene	13.89	228	729899	39.852	nq	99
84) Bis(2-ethylhexyl)phthalate	13.85	149	444507	40.086	nq	100
85) Di-n-octyl phthalate	14.46	149	887868	43.670	nq	100
87) Indeno(1,2,3-cd)pyrene	16.64	276	962681	41.643	nq	100
88) Benzo(b)fluoranthene	14.87	252	789521	37.897	nq	99
89) Benzo(k)fluoranthene	14.90	252	754157	38.982	nq	99
90) Benzo(a)pyrene	15.22	252	734649	39.069	nq	99
91) Dibenzo(a,h)anthracene	16.66	278	792888	40.028	nq	99
92) Benzo(g,h,i)perylene	17.06	276	827671	42.341	nq	99

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

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