

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF051820\
 Data File : BF120341.D
 Acq On : 18 May 2020 15:02
 Operator : MA/SJ
 Sample : PB128931BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB128931BS

Manual Integrations
 APPROVED

mohammad
 5/19/2020 3:05:07 PM

Quant Time: May 18 15:33:12 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF051420.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 18 15:23:07 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.60	152	335649	20.00	ng	0.00
21) Naphthalene-d8	7.89	136	1345530	20.00	ng	0.00
39) Acenaphthene-d10	9.65	164	679649	20.00	ng	0.00
64) Phenanthrene-d10	11.13	188	1094839	20.00	ng	0.00
76) Chrysene-d12	13.76	240	773913	20.00	ng	0.00
87) Perylene-d12	15.14	264	734390	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	2265878	131.35	ng	0.02
7) Phenol-d6	6.27	99	2912874	133.53	ng	0.00
23) Nitrobenzene-d5	7.17	82	1927580	98.08	ng	0.00
42) 2,4,6-Tribromophenol	10.44	330	709736	128.08	ng	0.00
45) 2-Fluorobiphenyl	8.97	172	3621787	109.48	ng	0.00
79) Terphenyl-d14	12.71	244	3180500	90.79	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	2.30	88	375574	54.730	ng	99
3) Pyridine	2.96	79	826285	38.305	ng	99
4) n-Nitrosodimethylamine	2.92	42	430981	51.025	ng	100
6) Aniline	6.27	93	1022054	36.634	ng	96
8) 2-Chlorophenol	6.39	128	959564	47.707	ng	99
9) Benzaldehyde	6.15	77	583488	43.388	ng	99
10) Phenol	6.28	94	1143244	44.893	ng	# 76
11) bis(2-Chloroethyl)ether	6.35	93	907548	44.614	ng	96
12) 1,3-Dichlorobenzene	6.54	146	964908	44.832	ng	100
13) 1,4-Dichlorobenzene	6.62	146	1004012	45.292	ng	99
14) 1,2-Dichlorobenzene	6.77	146	872843	43.725	ng	98
15) Benzyl Alcohol	6.76	79	715459	46.164	ng	100
16) 2,2'-oxybis(1-Chloropropan	6.89	45	1463288	45.646	ng	78
17) 2-Methylphenol	6.88	107	738684	43.504	ng	95
18) Hexachloroethane	7.11	117	370727	43.900	ng	99
19) n-Nitroso-di-n-propylamine	7.03	70	634376	44.387	ng	98
20) 3+4-Methylphenols	7.04	107	950189	43.041	ng	96
22) Acetophenone	7.02	105	1281570	46.493	ng	96
24) Nitrobenzene	7.19	77	980649	46.366	ng	97
25) Isophorone	7.44	82	1839997	44.998	ng	99
26) 2-Nitrophenol	7.51	139	536767	48.526	ng	95
27) 2,4-Dimethylphenol	7.56	122	819717	50.943	ng	97
28) bis(2-Chloroethoxy)methane	7.65	93	1191802	46.371	ng	99
29) 2,4-Dichlorophenol	7.76	162	772215	46.727	ng	98
30) 1,2,4-Trichlorobenzene	7.83	180	837367	45.724	ng	98
31) Naphthalene	7.91	128	2645819	46.592	ng	100
32) Benzoic acid	7.72	122	475763	45.108	ng	95
33) 4-Chloroaniline	7.97	127	763053	29.485	ng	100
34) Hexachlorobutadiene	8.03	225	467860	44.200	ng	95
35) Caprolactam	8.36	113	272430m	50.389	ng	
36) 4-Chloro-3-methylphenol	8.47	107	913171	51.079	ng	96
37) 2-Methylnaphthalene	8.60	142	1944586	50.039	ng	98
38) 1-Methylnaphthalene	8.70	142	1884562	50.001	ng	100
40) 1,2,4,5-Tetrachlorobenzene	8.77	216	871138	50.832	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.76	237	757329	104.881	nq	99
43) 2,4,6-Trichlorophenol	8.89	196	566752	49.268	nq	99
44) 2,4,5-Trichlorophenol	8.94	196	611293	46.272	nq	99
46) 1,1'-Biphenyl	9.07	154	2302727	51.710	nq	99
47) 2-Chloronaphthalene	9.09	162	1749650	48.981	nq	99
48) 2-Nitroaniline	9.20	65	625560	47.997	nq	98
49) Acenaphthylene	9.50	152	2731047	50.051	nq	100
50) Dimethylphthalate	9.38	163	2049868	47.712	nq	99
51) 2,6-Dinitrotoluene	9.45	165	462709	48.254	nq	87
52) Acenaphthene	9.67	154	1463196	44.737	nq	100
53) 3-Nitroaniline	9.61	138	319043	27.900	nq	97
54) 2,4-Dinitrophenol	9.73	184	399674	88.300	nq	94
55) Dibenzofuran	9.85	168	2148049	45.606	nq	99
56) 4-Nitrophenol	9.80	139	524687	85.394	nq	89
57) 2,4-Dinitrotoluene	9.85	165	494210	43.936	nq	97
58) Fluorene	10.19	166	1580394	44.059	nq	100
59) 2,3,4,6-Tetrachlorophenol	9.97	232	420017	42.247	nq	95
60) Diethylphthalate	10.08	149	1749465	42.178	nq	99
61) 4-Chlorophenyl-phenylether	10.19	204	765448	41.044	nq	96
62) 4-Nitroaniline	10.23	138	438210	41.154	nq	98
63) Azobenzene	10.35	77	1697593	43.304	nq	95
65) 4,6-Dinitro-2-methylphenol	10.26	198	278101	50.226	nq	97
66) n-Nitrosodiphenylamine	10.31	169	1491119	49.229	nq	99
67) 4-Bromophenyl-phenylether	10.67	248	462564	43.539	nq	97
68) Hexachlorobenzene	10.74	284	503981	45.344	nq	96
69) Atrazine	10.85	200	394297	43.504	nq	99
70) Pentachlorophenol	10.94	266	438390	92.739	nq	99
71) Phenanthrene	11.15	178	2190655	47.618	nq	99
72) Anthracene	11.20	178	2353519	47.787	nq	100
73) Carbazole	11.36	167	2138636	46.475	nq	100
74) Di-n-butylphthalate	11.69	149	2570082	46.342	nq	99
75) Fluoranthene	12.33	202	2454460	49.776	nq	99
77) Benzidine	12.46	184	1324692	58.616	nq	96
78) Pyrene	12.56	202	2482853	45.002	nq	100
80) Butylbenzylphthalate	13.19	149	1088769	41.908	nq	96
81) Benzo(a)anthracene	13.75	228	1866629	45.730	nq	99
82) 3,3'-Dichlorobenzidine	13.72	252	533307	34.670	nq	98
83) Chrysene	13.79	228	2005755	45.760	nq	100
84) Bis(2-ethylhexyl)phthalate	13.75	149	1366788	42.285	nq	99
85) Di-n-octyl phthalate	14.36	149	2547212	43.805	nq	99
86) Indeno(1,2,3-cd)pyrene	16.46	276	1710413	42.995	nq	99
88) Benzo(b)fluoranthene	14.76	252	1886948	47.048	nq	98
89) Benzo(k)fluoranthene	14.79	252	1718302	49.451	nq	100
90) Benzo(a)pyrene	15.09	252	1643231	45.685	nq	98
91) Dibenzo(a,h)anthracene	16.47	278	1417763	44.493	nq	98
92) Benzo(g,h,i)perylene	16.85	276	1453813	45.244	nq	98

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

