

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF032020\
 Data File : BF119472.D
 Acq On : 20 Mar 2020 15:13
 Operator : CG/JU
 Sample : PB127713BS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB127713BS

Manual Integrations
 APPROVED

mohammad
 3/23/2020 9:40:43 AM

Quant Time: Mar 23 04:26:31 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 18 15:11:18 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	311649	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	1220907	20.00	ng	0.00
39) Acenaphthene-d10	9.90	164	638706	20.00	ng	0.00
64) Phenanthrene-d10	11.39	188	1191082	20.00	ng	0.00
76) Chrysene-d12	14.03	240	621142	20.00	ng	0.00
87) Perylene-d12	15.50	264	670481	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.47	112	2320816	118.55	ng	0.02
7) Phenol-d6	6.48	99	2954832	118.15	ng	0.00
23) Nitrobenzene-d5	7.42	82	1860546	82.82	ng	0.00
42) 2,4,6-Tribromophenol	10.69	330	826163	125.38	ng	0.00
45) 2-Fluorobiphenyl	9.22	172	3299049	76.52	ng	0.00
79) Terphenyl-d14	12.98	244	3481385	109.81	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.66	88	322087	33.311	ng	98
3) Pyridine	3.40	79	869252	32.633	ng	98
4) n-Nitrosodimethylamine	3.36	42	559371	43.062	ng	99
6) Aniline	6.52	93	1353023	39.201	ng	99
8) 2-Chlorophenol	6.64	128	1045952	50.869	ng	97
9) Benzaldehyde	6.40	77	659034	41.541	ng	99
10) Phenol	6.49	94	1286001m	48.193	ng	
11) bis(2-Chloroethyl)ether	6.59	93	989825	46.379	ng	95
12) 1,3-Dichlorobenzene	6.79	146	1037429	46.618	ng	98
13) 1,4-Dichlorobenzene	6.87	146	1040310	46.736	ng	99
14) 1,2-Dichlorobenzene	7.03	146	971526	46.695	ng	99
15) Benzyl Alcohol	6.99	79	916060	48.696	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.13	45	1899457	44.124	ng	99
17) 2-Methylphenol	7.10	107	851084	45.177	ng	99
18) Hexachloroethane	7.37	117	396333	48.586	ng	99
19) n-Nitroso-di-n-propylamine	7.27	70	705699	46.816	ng	95
20) 3+4-Methylphenols	7.26	107	1010655	43.774	ng	# 88
22) Acetophenone	7.26	105	1319704	45.237	ng	# 90
24) Nitrobenzene	7.44	77	1080654	45.013	ng	96
25) Isophorone	7.68	82	2005911	44.018	ng	98
26) 2-Nitrophenol	7.75	139	546014	57.762	ng	97
27) 2,4-Dimethylphenol	7.79	122	961420	49.188	ng	98
28) bis(2-Chloroethoxy)methane	7.89	93	1249113	46.354	ng	99
29) 2,4-Dichlorophenol	7.99	162	841955	46.881	ng	99
30) 1,2,4-Trichlorobenzene	8.08	180	896323	45.128	ng	98
31) Naphthalene	8.16	128	2809016	44.709	ng	99
32) Benzoic acid	7.92	122	611442	48.631	ng	98
33) 4-Chloroaniline	8.20	127	860512	29.890	ng	100
34) Hexachlorobutadiene	8.28	225	501873	45.162	ng	97
35) Caprolactam	8.59	113	279475m	47.548	ng	
36) 4-Chloro-3-methylphenol	8.69	107	928232	47.289	ng	98
37) 2-Methylnaphthalene	8.86	142	1917030	46.491	ng	100
38) 1-Methylnaphthalene	8.96	142	1805613	46.263	ng	100
40) 1,2,4,5-Tetrachlorobenzene	9.02	216	813385	43.448	ng	99

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41) Hexachlorocyclopentadiene	9.01	237	878930	82.582	nq	100
43) 2,4,6-Trichlorophenol	9.13	196	626387	46.755	nq	99
44) 2,4,5-Trichlorophenol	9.17	196	655438	43.673	nq	97
46) 1,1'-Biphenyl	9.32	154	2237352	43.010	nq	99
47) 2-Chloronaphthalene	9.35	162	1784230	43.788	nq	99
48) 2-Nitroaniline	9.44	65	679573	48.318	nq	95
49) Acenaphthylene	9.76	152	2836251	44.325	nq	98
50) Dimethylphthalate	9.62	163	2113588	46.588	nq	98
51) 2,6-Dinitrotoluene	9.68	165	487266	50.108	nq	95
52) Acenaphthene	9.93	154	1869179	46.857	nq	100
53) 3-Nitroaniline	9.85	138	375376	30.576	nq	98
54) 2,4-Dinitrophenol	9.96	184	374447	88.173	nq	92
55) Dibenzofuran	10.11	168	2515388	43.980	nq	100
56) 4-Nitrophenol	10.01	139	888916	91.785	nq	89
57) 2,4-Dinitrotoluene	10.09	165	650033	53.064	nq	93
58) Fluorene	10.45	166	1888079	44.642	nq	98
59) 2,3,4,6-Tetrachlorophenol	10.22	232	551179	49.133	nq	99
60) Diethylphthalate	10.33	149	2153244	46.763	nq	98
61) 4-Chlorophenyl-phenylether	10.44	204	924596	44.704	nq	98
62) 4-Nitroaniline	10.47	138	512783	43.558	nq	98
63) Azobenzene	10.60	77	2074740	44.745	nq	96
65) 4,6-Dinitro-2-methylphenol	10.49	198	273806	54.821	nq	83
66) n-Nitrosodiphenylamine	10.56	169	1795856	45.928	nq	99
67) 4-Bromophenyl-phenylether	10.93	248	612119	45.125	nq	97
68) Hexachlorobenzene	11.00	284	640793	46.155	nq	98
69) Atrazine	11.09	200	552332	51.525	nq	99
70) Pentachlorophenol	11.19	266	708053	86.978	nq	98
71) Phenanthrene	11.42	178	2750644	44.537	nq	99
72) Anthracene	11.47	178	2825812	45.019	nq	100
73) Carbazole	11.62	167	2663225	42.865	nq	99
74) Di-n-butylphthalate	11.95	149	3298210	49.625	nq	98
75) Fluoranthene	12.60	202	2845755	42.576	nq	99
77) Benzidine	12.72	184	1906705	72.535	nq	99
78) Pyrene	12.83	202	2809038	55.105	nq	99
80) Butylbenzylphthalate	13.45	149	1340154	65.000	nq	96
81) Benzo(a)anthracene	14.02	228	1958563	48.489	nq	100
82) 3,3'-Dichlorobenzidine	13.98	252	705855	44.321	nq	98
83) Chrysene	14.06	228	1998854	47.950	nq	99
84) Bis(2-ethylhexyl)phthalate	14.02	149	1656904	67.414	nq	100
85) Di-n-octyl phthalate	14.63	149	2685288	62.123	nq	98
86) Indeno(1,2,3-cd)pyrene	16.98	276	2413257	61.468	nq	98
88) Benzo(b)fluoranthene	15.07	252	1988449	44.397	nq	97
89) Benzo(k)fluoranthene	15.10	252	1857503	43.555	nq	99
90) Benzo(a)pyrene	15.44	252	1814041	44.568	nq	98
91) Dibenzo(a,h)anthracene	17.00	278	1946279	54.669	nq	99
92) Benzo(g,h,i)perylene	17.43	276	2082199	58.217	nq	98

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

