

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF040820\
 Data File : BF119845.D
 Acq On : 9 Apr 2020 1:11
 Operator : MA/SJ
 Sample : PB128073BS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB128073BS

Manual Integrations
 APPROVED

mohammad
 4/10/2020 9:57:18 AM

Quant Time: Apr 09 02:43:22 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF040820.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 09 01:21:19 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	143453	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	521144	20.00	ng	0.00
39) Acenaphthene-d10	9.81	164	276618	20.00	ng	0.00
64) Phenanthrene-d10	11.30	188	621758	20.00	ng	0.00
76) Chrysene-d12	13.94	240	464627	20.00	ng	0.00
87) Perylene-d12	15.37	264	428529	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.39	112	1070254	128.93	ng	0.01
7) Phenol-d6	6.40	99	1334139	124.73	ng	0.00
23) Nitrobenzene-d5	7.33	82	816896	81.09	ng	0.00
42) 2,4,6-Tribromophenol	10.60	330	447764	132.21	ng	0.00
45) 2-Fluorobiphenyl	9.13	172	1757821	90.22	ng	0.00
79) Terphenyl-d14	12.89	244	2191242	79.36	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.54	88	144134	38.985	ng	97
3) Pyridine	3.25	79	393126	38.755	ng	96
4) n-Nitrosodimethylamine	3.20	42	232447	44.850	ng	98
6) Aniline	6.43	93	432485	30.807	ng	99
8) 2-Chlorophenol	6.55	128	443408	47.808	ng	97
9) Benzaldehyde	6.31	77	175503	26.179	ng	97
10) Phenol	6.42	94	538554	44.382	ng	96
11) bis(2-Chloroethyl)ether	6.50	93	382467	40.821	ng	99
12) 1,3-Dichlorobenzene	6.70	146	443047	43.633	ng	99
13) 1,4-Dichlorobenzene	6.78	146	447273	43.242	ng	99
14) 1,2-Dichlorobenzene	6.94	146	410675	43.082	ng	97
15) Benzyl Alcohol	6.91	79	350895m	42.238	ng	
16) 2,2'-oxybis(1-Chloropropan	7.05	45	714713	39.201	ng	99
17) 2-Methylphenol	7.03	107	322635	38.980	ng	96
18) Hexachloroethane	7.28	117	159720	41.319	ng	92
19) n-Nitroso-di-n-propylamine	7.19	70	295138	42.910	ng	98
20) 3+4-Methylphenols	7.17	107	432698	42.744	ng	94
22) Acetophenone	7.17	105	570355	46.737	ng	96
24) Nitrobenzene	7.35	77	445205	43.933	ng	100
25) Isophorone	7.59	82	797619	41.074	ng	100
26) 2-Nitrophenol	7.67	139	233686	46.158	ng	94
27) 2,4-Dimethylphenol	7.71	122	377773	49.475	ng	97
28) bis(2-Chloroethoxy)methane	7.80	93	536795	46.976	ng	99
29) 2,4-Dichlorophenol	7.91	162	367198	46.916	ng	97
30) 1,2,4-Trichlorobenzene	7.99	180	369900	41.710	ng	99
31) Naphthalene	8.07	128	1180000	43.036	ng	99
32) Benzoic acid	7.83	122	307591	45.868	ng	97
33) 4-Chloroaniline	8.12	127	292136	24.474	ng	98
34) Hexachlorobutadiene	8.19	225	215224	40.542	ng	98
35) Caprolactam	8.49	113	108716m	45.409	ng	
36) 4-Chloro-3-methylphenol	8.61	107	398826	47.066	ng	98
37) 2-Methylnaphthalene	8.77	142	864128	46.486	ng	99
38) 1-Methylnaphthalene	8.87	142	775183	44.243	ng	100
40) 1,2,4,5-Tetrachlorobenzene	8.93	216	380709	43.575	ng	99

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41) Hexachlorocyclopentadiene	8.92	237	481502	96.919	nq	97
43) 2,4,6-Trichlorophenol	9.05	196	259841	43.069	nq	100
44) 2,4,5-Trichlorophenol	9.09	196	292929	43.109	nq	97
46) 1,1'-Biphenyl	9.23	154	1070753	45.860	nq	99
47) 2-Chloronaphthalene	9.26	162	819541	45.565	nq	98
48) 2-Nitroaniline	9.35	65	302050	46.341	nq	95
49) Acenaphthylene	9.67	152	1314860	48.069	nq	99
50) Dimethylphthalate	9.54	163	998353	46.661	nq	99
51) 2,6-Dinitrotoluene	9.60	165	222030	47.388	nq	88
52) Acenaphthene	9.85	154	888597m	48.699	nq	
53) 3-Nitroaniline	9.76	138	149679	27.971	nq	99
54) 2,4-Dinitrophenol	9.87	184	264524	94.300	nq	92
55) Dibenzofuran	10.02	168	1204249	48.095	nq	98
56) 4-Nitrophenol	9.93	139	389334	97.785	nq	98
57) 2,4-Dinitrotoluene	10.00	165	301216	48.795	nq	98
58) Fluorene	10.36	166	911336	45.510	nq	99
59) 2,3,4,6-Tetrachlorophenol	10.13	232	266936	51.363	nq	96
60) Diethylphthalate	10.24	149	1019171	45.959	nq	99
61) 4-Chlorophenyl-phenylether	10.35	204	442074	43.073	nq	95
62) 4-Nitroaniline	10.39	138	231454	45.386	nq	95
63) Azobenzene	10.52	77	946662	47.635	nq	99
65) 4,6-Dinitro-2-methylphenol	10.41	198	167795	40.964	nq	77
66) n-Nitrosodiphenylamine	10.47	169	847401	40.981	nq	99
67) 4-Bromophenyl-phenylether	10.85	248	297281	38.447	nq	95
68) Hexachlorobenzene	10.91	284	329632	42.023	nq	96
69) Atrazine	11.00	200	266889	46.389	nq	98
70) Pentachlorophenol	11.10	266	380099	84.087	nq	99
71) Phenanthrene	11.32	178	1455799	43.994	nq	98
72) Anthracene	11.37	178	1485175	43.935	nq	100
73) Carbazole	11.53	167	1316964	41.197	nq	99
74) Di-n-butylphthalate	11.86	149	1685408	40.105	nq	99
75) Fluoranthene	12.51	202	1555117	43.555	nq	98
77) Benzidine	12.63	184	628798	40.036	nq	96
78) Pyrene	12.74	202	1563686	39.922	nq	99
80) Butylbenzylphthalate	13.36	149	745564	39.228	nq	99
81) Benzo(a)anthracene	13.93	228	1258104	41.160	nq	100
82) 3,3'-Dichlorobenzidine	13.89	252	345727	30.314	nq	96
83) Chrysene	13.97	228	1253813	40.370	nq	99
84) Bis(2-ethylhexyl)phthalate	13.93	149	1004633	39.033	nq	100
85) Di-n-octyl phthalate	14.54	149	1794261	40.943	nq	99
86) Indeno(1,2,3-cd)pyrene	16.79	276	1174234	42.891	nq	99
88) Benzo(b)fluoranthene	14.96	252	1499974	48.657	nq	99
89) Benzo(k)fluoranthene	14.99	252	1212135	45.572	nq	99
90) Benzo(a)pyrene	15.32	252	1141817	44.052	nq	99
91) Dibenzo(a,h)anthracene	16.81	278	989324	43.090	nq	99
92) Benzo(g,h,i)perylene	17.22	276	915131	40.380	nq	99

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

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