

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091919\
 Data File : BF117172.D
 Acq On : 19 Sep 2019 13:54
 Operator : HP/JU
 Sample : SSTDC0040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDC0040

Manual Integrations
 APPROVED

mohammad
 9/20/2019 1:06:42 PM

Quant Time: Sep 20 01:45:06 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 13 16:23:17 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.82	152	69203	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	295603	20.00	ng	0.00
39) Acenaphthene-d10	9.85	164	158712	20.00	ng	0.00
64) Phenanthrene-d10	11.33	188	233302	20.00	ng	0.00
76) Chrysene-d12	13.97	240	207872	20.00	ng	0.00
87) Perylene-d12	15.42	264	160175	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.45	112	384737	86.80	ng	0.00
7) Phenol-d6	6.46	99	427758	74.67	ng	0.00
23) Nitrobenzene-d5	7.39	82	456283	81.10	ng	0.00
42) 2,4,6-Tribromophenol	10.65	330	102368	77.17	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	792633	78.09	ng	0.00
79) Terphenyl-d14	12.92	244	871596	78.38	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.56	88	44306	23.887	ng	96
3) Pyridine	3.31	79	123262m	24.280	ng	
4) n-Nitrosodimethylamine	3.26	42	44788m	25.708	ng	
6) Aniline	6.49	93	281367	38.149	ng	92
8) 2-Chlorophenol	6.62	128	180898	37.830	ng	96
9) Benzaldehyde	6.37	77	122189	43.520	ng	96
10) Phenol	6.48	94	236773	37.493	ng	84
11) bis(2-Chloroethyl)ether	6.56	93	182036	39.220	ng	98
12) 1,3-Dichlorobenzene	6.76	146	204860	38.178	ng	99
13) 1,4-Dichlorobenzene	6.84	146	213188	38.932	ng	98
14) 1,2-Dichlorobenzene	7.00	146	195706	38.386	ng	99
15) Benzyl Alcohol	6.97	79	172207	39.202	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.10	45	176105	38.404	ng	87
17) 2-Methylphenol	7.09	107	152752	38.115	ng	94
18) Hexachloroethane	7.33	117	89049	42.271	ng	95
19) n-Nitroso-di-n-propylamine	7.24	70	147224	42.206	ng	95
20) 3+4-Methylphenols	7.24	107	202204	40.506	ng	91
22) Acetophenone	7.23	105	302226	40.241	ng	# 98
24) Nitrobenzene	7.41	77	224469	40.402	ng	98
25) Isophorone	7.65	82	362528	36.394	ng	99
26) 2-Nitrophenol	7.72	139	95449	39.996	ng	93
27) 2,4-Dimethylphenol	7.76	122	147648	39.015	ng	99
28) bis(2-Chloroethoxy)methane	7.85	93	227816	37.120	ng	100
29) 2,4-Dichlorophenol	7.97	162	132942	33.526	ng	97
30) 1,2,4-Trichlorobenzene	8.05	180	154097	35.970	ng	97
31) Naphthalene	8.13	128	554626	39.012	ng	99
32) Benzoic acid	7.90	122	77308	29.931	ng	94
33) 4-Chloroaniline	8.17	127	244187	38.726	ng	98
34) Hexachlorobutadiene	8.25	225	102619	42.220	ng	97
35) Caprolactam	8.55	113	47905	35.692	ng	92
36) 4-Chloro-3-methylphenol	8.66	107	168589	36.462	ng	97
37) 2-Methylnaphthalene	8.82	142	350399	36.601	ng	99
38) 1-Methylnaphthalene	8.91	142	311339	34.723	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.98	216	146068	35.641	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.97	237	50700	33.088	nq	98
43) 2,4,6-Trichlorophenol	9.09	196	106141	38.163	nq	99
44) 2,4,5-Trichlorophenol	9.14	196	111698	37.589	nq	98
46) 1,1'-Biphenyl	9.27	154	488989	39.278	nq	99
47) 2-Chloronaphthalene	9.30	162	387513	39.441	nq	99
48) 2-Nitroaniline	9.40	65	127554	40.475	nq	91
49) Acenaphthylene	9.72	152	559222	37.994	nq	100
50) Dimethylphthalate	9.57	163	436359	38.830	nq	99
51) 2,6-Dinitrotoluene	9.64	165	88555	38.692	nq	# 83
52) Acenaphthene	9.89	154	318055	36.453	nq	99
53) 3-Nitroaniline	9.82	138	107390	37.189	nq	94
54) 2,4-Dinitrophenol	9.93	184	28419	34.628	nq	90
55) Dibenzofuran	10.06	168	487373	37.860	nq	97
56) 4-Nitrophenol	9.98	139	69233	36.992	nq	96
57) 2,4-Dinitrotoluene	10.05	165	117962	39.706	nq	92
58) Fluorene	10.40	166	410721	39.608	nq	100
59) 2,3,4,6-Tetrachlorophenol	10.18	232	86923m	38.225	nq	
60) Diethylphthalate	10.27	149	449951	40.698	nq	99
61) 4-Chlorophenyl-phenylether	10.39	204	185625	41.373	nq	96
62) 4-Nitroaniline	10.43	138	104685	34.859	nq	94
63) Azobenzene	10.55	77	424436	37.588	nq	97
65) 4,6-Dinitro-2-methylphenol	10.46	198	39531	38.742	nq	# 71
66) n-Nitrosodiphenylamine	10.51	169	317908	39.456	nq	99
67) 4-Bromophenyl-phenylether	10.87	248	94746	39.766	nq	89
68) Hexachlorobenzene	10.95	284	101826	39.074	nq	# 93
69) Atrazine	11.03	200	78449	37.384	nq	99
70) Pentachlorophenol	11.15	266	53315	43.646	nq	96
71) Phenanthrene	11.36	178	517923	39.084	nq	99
72) Anthracene	11.41	178	510486	37.733	nq	98
73) Carbazole	11.56	167	493954	38.465	nq	99
74) Di-n-butylphthalate	11.89	149	707062	46.278	nq	99
75) Fluoranthene	12.55	202	637147	46.794	nq	99
77) Benzidine	12.66	184	219794	32.824	nq	99
78) Pyrene	12.77	202	638467	36.605	nq	100
80) Butylbenzylphthalate	13.38	149	348818	42.324	nq	99
81) Benzo(a)anthracene	13.96	228	520836	37.502	nq	99
82) 3,3'-Dichlorobenzidine	13.92	252	169929	37.970	nq	99
83) Chrysene	14.00	228	498471	36.800	nq	99
84) Bis(2-ethylhexyl)phthalate	13.94	149	495719	46.650	nq	100
85) Di-n-octyl phthalate	14.55	149	786050	46.998	nq	100
86) Indeno(1,2,3-cd)pyrene	16.86	276	302606m	30.621	nq	
88) Benzo(b)fluoranthene	15.00	252	377136	38.870	nq	99
89) Benzo(k)fluoranthene	15.03	252	377215	41.043	nq	99
90) Benzo(a)pyrene	15.36	252	329316	38.679	nq	99
91) Dibenzo(a,h)anthracene	16.87	278	250849	36.985	nq	99
92) Benzo(g,h,i)perylene	17.30	276	233936	34.612	nq	97

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

