

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062519\
 Data File : BF115186.D
 Acq On : 25 Jun 2019 14:41
 Operator : HP/JU
 Sample : PB120785BS
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB120785BS

Manual Integrations
 APPROVED

mohammad
 6/26/2019 2:03:21 PM

Quant Time: Jun 26 06:47:53 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF062119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 24 04:21:29 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.61	152	279456	20.00	ng	0.00
21) Naphthalene-d8	7.90	136	1113435	20.00	ng	0.00
39) Acenaphthene-d10	9.64	164	548925	20.00	ng	0.00
64) Phenanthrene-d10	11.11	188	1075730	20.00	ng	0.00
76) Chrysene-d12	13.74	240	727448	20.00	ng	-0.01
87) Perylene-d12	15.10	264	761941	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.22	112	1975813	109.11	ng	0.02
7) Phenol-d6	6.26	99	2434266	110.75	ng	0.00
23) Nitrobenzene-d5	7.18	82	1502044	82.99	ng	0.00
42) 2,4,6-Tribromophenol	10.42	330	701301	121.15	ng	0.00
45) 2-Fluorobiphenyl	8.97	172	2779464	86.01	ng	-0.01
79) Terphenyl-d14	12.69	244	2901092	84.79	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.30	88	303923	35.561	ng	98
3) Pyridine	2.95	79	831026	34.499	ng	96
4) n-Nitrosodimethylamine	2.90	42	345262	36.561	ng	100
6) Aniline	6.27	93	726423	24.047	ng	# 1
8) 2-Chlorophenol	6.40	128	835893	41.050	ng	98
9) Benzaldehyde	6.15	77	115837	8.078	ng	99
10) Phenol	6.27	94	944348	36.678	ng	93
11) bis(2-Chloroethyl)ether	6.35	93	752686	39.659	ng	98
12) 1,3-Dichlorobenzene	6.55	146	877061	38.810	ng	100
13) 1,4-Dichlorobenzene	6.63	146	885577	39.078	ng	99
14) 1,2-Dichlorobenzene	6.78	146	822349	39.185	ng	99
15) Benzyl Alcohol	6.76	79	540390	33.763	ng	99
16) 2,2'-oxybis(1-Chloropropan	6.90	45	1073574	39.001	ng	96
17) 2-Methylphenol	6.88	107	722990	43.015	ng	98
18) Hexachloroethane	7.12	117	318168	38.944	ng	97
19) n-Nitroso-di-n-propylamine	7.04	70	531136	37.744	ng	99
20) 3+4-Methylphenols	7.03	107	821120	42.309	ng	91
22) Acetophenone	7.02	105	1099037	43.116	ng	# 97
24) Nitrobenzene	7.20	77	836197	41.935	ng	98
25) Isophorone	7.44	82	1544457	41.653	ng	99
26) 2-Nitrophenol	7.51	139	477367	48.476	ng	99
27) 2,4-Dimethylphenol	7.56	122	797749	49.478	ng	98
28) bis(2-Chloroethoxy)methane	7.65	93	965089	41.181	ng	99
29) 2,4-Dichlorophenol	7.75	162	738970	44.412	ng	97
30) 1,2,4-Trichlorobenzene	7.84	180	782503	42.576	ng	99
31) Naphthalene	7.92	128	2337476	42.767	ng	100
32) Benzoic acid	7.70	122	597582	51.923	ng	100
33) 4-Chloroaniline	7.97	127	573019	22.940	ng	98
34) Hexachlorobutadiene	8.04	225	421033	41.803	ng	98
35) Caprolactam	8.34	113	244966m	43.789	ng	
36) 4-Chloro-3-methylphenol	8.46	107	750982	44.567	ng	96
37) 2-Methylnaphthalene	8.61	142	1602833	43.494	ng	99
38) 1-Methylnaphthalene	8.71	142	1525706	43.349	ng	100
40) 1,2,4,5-Tetrachlorobenzene	8.78	216	668888	43.122	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.77	237	706307	76.790	nq	99
43) 2,4,6-Trichlorophenol	8.88	196	511280	42.694	nq	99
44) 2,4,5-Trichlorophenol	8.93	196	591765	47.984	nq	99
46) 1,1'-Biphenyl	9.07	154	1832402	41.720	nq	98
47) 2-Chloronaphthalene	9.09	162	1503380	42.197	nq	97
48) 2-Nitroaniline	9.18	65	454551	43.762	nq	98
49) Acenaphthylene	9.50	152	2325568	41.896	nq	99
50) Dimethylphthalate	9.38	163	1706285	42.250	nq	100
51) 2,6-Dinitrotoluene	9.43	165	414360	44.441	nq	98
52) Acenaphthene	9.68	154	1438667	41.419	nq	98
53) 3-Nitroaniline	9.60	138	314558	27.646	nq	93
54) 2,4-Dinitrophenol	9.70	184	453070	93.109	nq	94
55) Dibenzofuran	9.85	168	2022904	40.577	nq	98
56) 4-Nitrophenol	9.77	139	715368	78.424	nq	94
57) 2,4-Dinitrotoluene	9.83	165	549431	45.638	nq	93
58) Fluorene	10.19	166	1403112	41.805	nq	98
59) 2,3,4,6-Tetrachlorophenol	9.97	232	495036	47.871	nq	97
60) Diethylphthalate	10.08	149	1677432	41.320	nq	99
61) 4-Chlorophenyl-phenylether	10.18	204	681595	42.893	nq	97
62) 4-Nitroaniline	10.21	138	466388	40.860	nq	98
63) Azobenzene	10.34	77	1561238	41.174	nq	98
65) 4,6-Dinitro-2-methylphenol	10.24	198	271724	47.527	nq	89
66) n-Nitrosodiphenylamine	10.30	169	1422598	42.448	nq	100
67) 4-Bromophenyl-phenylether	10.67	248	490295	42.038	nq	92
68) Hexachlorobenzene	10.73	284	541679	42.788	nq	95
69) Atrazine	10.83	200	458504	42.529	nq	99
70) Pentachlorophenol	10.92	266	674779	83.667	nq	96
71) Phenanthrene	11.14	178	2294379	39.614	nq	100
72) Anthracene	11.19	178	2343554	40.085	nq	98
73) Carbazole	11.34	167	2155717	41.292	nq	98
74) Di-n-butylphthalate	11.69	149	2466828	40.890	nq	98
75) Fluoranthene	12.32	202	2363573	40.901	nq	98
77) Benzidine	12.44	184	805149	24.926	nq	99
78) Pyrene	12.54	202	2393960	42.822	nq	98
80) Butylbenzylphthalate	13.18	149	1114478	45.173	nq	91
81) Benzo(a)anthracene	13.72	228	2221933	42.568	nq	99
82) 3,3'-Dichlorobenzidine	13.69	252	697877	37.024	nq	99
83) Chrysene	13.76	228	2101713	43.957	nq	99
84) Bis(2-ethylhexyl)phthalate	13.74	149	1417499	43.848	nq	99
85) Di-n-octyl phthalate	14.35	149	2451146	45.128	nq	99
86) Indeno(1,2,3-cd)pyrene	16.37	276	2343359	41.419	nq	99
88) Benzo(b)fluoranthene	14.73	252	2416891	50.836	nq	99
89) Benzo(k)fluoranthene	14.76	252	1993147	46.748	nq	98
90) Benzo(a)pyrene	15.05	252	2169226	50.622	nq	99
91) Dibenzo(a,h)anthracene	16.39	278	1936394	48.405	nq	98
92) Benzo(g,h,i)perylene	16.76	276	1950212	48.167	nq	99

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

