

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092519\
 Data File : BF117238.D
 Acq On : 25 Sep 2019 14:02
 Operator : JU
 Sample : K5008-07MS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-7MS

Manual Integrations
 APPROVED

mohammad
 9/26/2019 3:53:35 PM

Quant Time: Sep 26 03:13:43 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	37738	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	154280	20.00	ng	0.00
39) Acenaphthene-d10	9.85	164	80441	20.00	ng	0.00
64) Phenanthrene-d10	11.34	188	122492	20.00	ng	0.00
76) Chrysene-d12	13.98	240	70001	20.00	ng	0.00
87) Perylene-d12	15.44	264	65907	20.00	ng	0.02

System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	223213	92.34	ng	0.01
7) Phenol-d6	6.46	99	312991	100.19	ng	0.00
23) Nitrobenzene-d5	7.38	82	193023	65.74	ng	0.00
42) 2,4,6-Tribromophenol	10.65	330	57152	85.01	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	314875	61.20	ng	0.00
79) Terphenyl-d14	12.92	244	250908	67.00	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.65	88	34428	34.038	ng	# 96
3) Pyridine	3.40	79	80756	29.170	ng	95
4) n-Nitrosodimethylamine	3.35	42	40336	42.456	ng	83
6) Aniline	6.48	93	99003	24.616	ng	# 59
8) 2-Chlorophenol	6.61	128	108262	41.517	ng	95
9) Benzaldehyde	6.36	77	74025	49.719	ng	96
10) Phenol	6.47	94	144535	41.970	ng	91
11) bis(2-Chloroethyl)ether	6.55	93	100760	39.810	ng	98
12) 1,3-Dichlorobenzene	6.76	146	112610	38.484	ng	97
13) 1,4-Dichlorobenzene	6.83	146	114400	38.310	ng	98
14) 1,2-Dichlorobenzene	6.99	146	108913	39.174	ng	98
15) Benzyl Alcohol	6.96	79	99908	41.706	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.09	45	94300	37.711	ng	# 42
17) 2-Methylphenol	7.08	107	89980	41.172	ng	# 90
18) Hexachloroethane	7.32	117	44138	38.422	ng	99
19) n-Nitroso-di-n-propylamine	7.23	70	79342	41.711	ng	93
20) 3+4-Methylphenols	7.23	107	113330	41.632	ng	# 72
22) Acetophenone	7.23	105	163824	41.794	ng	98
24) Nitrobenzene	7.40	77	118503	40.867	ng	99
25) Isophorone	7.64	82	218470	42.022	ng	99
26) 2-Nitrophenol	7.72	139	54861	44.047	ng	93
27) 2,4-Dimethylphenol	7.76	122	91308	46.228	ng	99
28) bis(2-Chloroethoxy)methane	7.85	93	128262	40.043	ng	98
29) 2,4-Dichlorophenol	7.96	162	86188	41.645	ng	99
30) 1,2,4-Trichlorobenzene	8.04	180	86603	38.732	ng	99
31) Naphthalene	8.12	128	308330	41.554	ng	99
32) Benzoic acid	7.89	122	30339	24.142	ng	93
33) 4-Chloroaniline	8.17	127	43342	13.170	ng	98
34) Hexachlorobutadiene	8.24	225	51178	40.344	ng	97
35) Caprolactam	8.55	113	26070m	37.216	ng	
36) 4-Chloro-3-methylphenol	8.66	107	100640	41.704	ng	95
37) 2-Methylnaphthalene	8.81	142	212330	42.496	ng	100
38) 1-Methylnaphthalene	8.91	142	195674	41.814	ng	98
40) 1,2,4,5-Tetrachlorobenzene	8.98	216	85215	41.024	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.96	237	56652	63.496	ng	99
43) 2,4,6-Trichlorophenol	9.09	196	56173	39.849	ng	98
44) 2,4,5-Trichlorophenol	9.15	196	60613	40.245	ng	97
46) 1,1'-Biphenyl	9.27	154	252701	40.049	ng	96
47) 2-Chloronaphthalene	9.30	162	194269	39.012	ng	99
48) 2-Nitroaniline	9.40	65	60679	37.989	ng	94
49) Acenaphthylene	9.72	152	307737	41.252	ng	99
50) Dimethylphthalate	9.57	163	269728	47.357	ng	98
51) 2,6-Dinitrotoluene	9.64	165	49689	42.835	ng #	86
52) Acenaphthene	9.89	154	177820	40.211	ng	99
53) 3-Nitroaniline	9.81	138	32073	21.914	ng	94
54) 2,4-Dinitrophenol	9.93	184	17826	40.678	ng	93
55) Dibenzofuran	10.06	168	271834	41.663	ng	99
56) 4-Nitrophenol	9.99	139	51407	54.194	ng	94
57) 2,4-Dinitrotoluene	10.05	165	64572	42.883	ng #	88
58) Fluorene	10.40	166	219679	41.798	ng	98
59) 2,3,4,6-Tetrachlorophenol	10.19	232	47826	41.497	ng	97
60) Diethylphthalate	10.27	149	229650	40.983	ng	100
61) 4-Chlorophenyl-phenylether	10.39	204	95477	41.987	ng	97
62) 4-Nitroaniline	10.43	138	48774	32.045	ng	89
63) Azobenzene	10.55	77	238896	41.743	ng	98
65) 4,6-Dinitro-2-methylphenol	10.46	198	19426	36.703	ng	88
66) n-Nitrosodiphenylamine	10.51	169	188263	44.503	ng	98
67) 4-Bromophenyl-phenylether	10.88	248	54073	43.226	ng	96
68) Hexachlorobenzene	10.96	284	57312	41.888	ng	99
69) Atrazine	11.04	200	48733	46.798	ng	98
70) Pentachlorophenol	11.15	266	47018	73.311	ng	98
71) Phenanthrene	11.36	178	291846	41.947	ng	99
72) Anthracene	11.42	178	299466	42.160	ng	99
73) Carbazole	11.57	167	252700	37.480	ng	99
74) Di-n-butylphthalate	11.89	149	349827	43.609	ng	98
75) Fluoranthene	12.55	202	271642	37.998	ng	99
77) Benzidine	12.67	184	60866	26.993	ng	99
78) Pyrene	12.78	202	263385	44.842	ng	99
80) Butylbenzylphthalate	13.39	149	117429	42.311	ng	97
81) Benzo(a)anthracene	13.97	228	184201	39.386	ng	97
82) 3,3'-Dichlorobenzidine	13.93	252	43037	28.556	ng	95
83) Chrysene	14.01	228	178719	39.180	ng	100
84) Bis(2-ethylhexyl)phthalate	13.95	149	169176	47.276	ng	99
85) Di-n-octyl phthalate	14.57	149	263022	46.699	ng	100
86) Indeno(1,2,3-cd)pyrene	16.91	276	162581	48.855	ng	98
88) Benzo(b)fluoranthene	15.03	252	151556	37.963	ng	98
89) Benzo(k)fluoranthene	15.06	252	153777	40.663	ng	96
90) Benzo(a)pyrene	15.39	252	147457	42.092	ng	97
91) Dibenzo(a,h)anthracene	16.92	278	136870	49.043	ng	98
92) Benzo(g,h,i)perylene	17.34	276	127789	45.951	ng	97

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

