

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092519\
 Data File : BF117239.D
 Acq On : 25 Sep 2019 14:51
 Operator : JU
 Sample : K5008-07MSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-7MSD

Manual Integrations
 APPROVED

mohammad
 9/26/2019 3:53:38 PM

Quant Time: Sep 26 03:19:02 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	39194	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	161296	20.00	ng	0.00
39) Acenaphthene-d10	9.85	164	82643	20.00	ng	0.00
64) Phenanthrene-d10	11.34	188	126776	20.00	ng	0.00
76) Chrysene-d12	13.99	240	76110	20.00	ng	0.01
87) Perylene-d12	15.47	264	67110	20.00	ng	0.05

System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	226443	90.20	ng	0.01
7) Phenol-d6	6.46	99	319205	98.39	ng	0.00
23) Nitrobenzene-d5	7.38	82	196952	64.16	ng	0.00
42) 2,4,6-Tribromophenol	10.65	330	60372	87.41	ng	0.00
45) 2-Fluorobiphenyl	9.17	172	322780	61.07	ng	0.00
79) Terphenyl-d14	12.92	244	267726	65.75	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.65	88	34578	32.916	ng	# 93
3) Pyridine	3.40	79	80572	28.022	ng	96
4) n-Nitrosodimethylamine	3.35	42	41194	41.748	ng	85
6) Aniline	6.48	93	101050	24.191	ng	# 57
8) 2-Chlorophenol	6.61	128	110150	40.672	ng	95
9) Benzaldehyde	6.36	77	75998	49.006	ng	94
10) Phenol	6.47	94	149450	41.785	ng	92
11) bis(2-Chloroethyl)ether	6.55	93	102939	39.160	ng	98
12) 1,3-Dichlorobenzene	6.76	146	114710	37.746	ng	99
13) 1,4-Dichlorobenzene	6.83	146	116487	37.560	ng	99
14) 1,2-Dichlorobenzene	6.99	146	108671	37.635	ng	98
15) Benzyl Alcohol	6.96	79	104213	41.887	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.09	45	96323	37.089	ng	# 40
17) 2-Methylphenol	7.08	107	93506	41.196	ng	# 91
18) Hexachloroethane	7.33	117	44550	37.340	ng	93
19) n-Nitroso-di-n-propylamine	7.23	70	82520	41.770	ng	92
20) 3+4-Methylphenols	7.23	107	119957	42.429	ng	# 70
22) Acetophenone	7.23	105	167376	40.843	ng	99
24) Nitrobenzene	7.40	77	126112	41.600	ng	97
25) Isophorone	7.64	82	225580	41.502	ng	99
26) 2-Nitrophenol	7.72	139	56735	43.570	ng	99
27) 2,4-Dimethylphenol	7.76	122	94170	45.604	ng	97
28) bis(2-Chloroethoxy)methane	7.85	93	131879	39.381	ng	98
29) 2,4-Dichlorophenol	7.96	162	89866	41.533	ng	95
30) 1,2,4-Trichlorobenzene	8.04	180	89995	38.499	ng	99
31) Naphthalene	8.12	128	315768	40.706	ng	99
32) Benzoic acid	7.89	122	39254	28.311	ng	97
33) 4-Chloroaniline	8.17	127	49057	14.258	ng	95
34) Hexachlorobutadiene	8.23	225	52699	39.736	ng	98
35) Caprolactam	8.55	113	27636m	37.735	ng	
36) 4-Chloro-3-methylphenol	8.67	107	103344	40.962	ng	99
37) 2-Methylnaphthalene	8.81	142	216882	41.519	ng	99
38) 1-Methylnaphthalene	8.91	142	201825	41.252	ng	99
40) 1,2,4,5-Tetrachlorobenzene	8.98	216	85981	40.290	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
41) Hexachlorocyclopentadiene	8.96	237	55313	60.733	nq	96
43) 2,4,6-Trichlorophenol	9.10	196	57732	39.864	nq	96
44) 2,4,5-Trichlorophenol	9.15	196	62424	40.343	nq	95
46) 1,1'-Biphenyl	9.27	154	260464	40.179	nq	97
47) 2-Chloronaphthalene	9.30	162	201979	39.479	nq	98
48) 2-Nitroaniline	9.40	65	64231	39.142	nq	92
49) Acenaphthylene	9.72	152	310727	40.543	nq	100
50) Dimethylphthalate	9.57	163	280512	47.938	nq	98
51) 2,6-Dinitrotoluene	9.64	165	49739	41.735	nq #	81
52) Acenaphthene	9.89	154	182664	40.206	nq	99
53) 3-Nitroaniline	9.81	138	33024	21.963	nq	86
54) 2,4-Dinitrophenol	9.93	184	20129	43.801	nq	94
55) Dibenzofuran	10.06	168	280307	41.817	nq	98
56) 4-Nitrophenol	9.99	139	55255	56.699	nq	91
57) 2,4-Dinitrotoluene	10.05	165	66688	43.108	nq #	88
58) Fluorene	10.40	166	229606	42.522	nq	98
59) 2,3,4,6-Tetrachlorophenol	10.19	232	38785m	32.755	nq	
60) Diethylphthalate	10.27	149	238212	41.379	nq	97
61) 4-Chlorophenyl-phenylether	10.39	204	98738	42.264	nq	99
62) 4-Nitroaniline	10.43	138	51857	33.162	nq	89
63) Azobenzene	10.55	77	246964	42.003	nq	98
65) 4,6-Dinitro-2-methylphenol	10.46	198	19917	36.424	nq	78
66) n-Nitrosodiphenylamine	10.51	169	191764	43.799	nq	98
67) 4-Bromophenyl-phenylether	10.88	248	57150	44.142	nq	99
68) Hexachlorobenzene	10.96	284	60354	42.621	nq	96
69) Atrazine	11.04	200	51640	48.363	nq	98
70) Pentachlorophenol	11.16	266	52031	78.386	nq	98
71) Phenanthrene	11.36	178	305737	42.459	nq	99
72) Anthracene	11.42	178	318758	43.360	nq	99
73) Carbazole	11.57	167	268540	38.483	nq	99
74) Di-n-butylphthalate	11.89	149	375901	45.276	nq	99
75) Fluoranthene	12.55	202	284954	38.513	nq	100
77) Benzidine	12.67	184	68832	28.075	nq	98
78) Pyrene	12.78	202	288002	45.098	nq	99
80) Butylbenzylphthalate	13.39	149	130670	43.303	nq	95
81) Benzo(a)anthracene	13.98	228	201122	39.552	nq	96
82) 3,3'-Dichlorobenzidine	13.93	252	45370	27.688	nq	98
83) Chrysene	14.02	228	194755	39.269	nq	97
84) Bis(2-ethylhexyl)phthalate	13.96	149	183110	47.063	nq	100
85) Di-n-octyl phthalate	14.59	149	276487	45.150	nq	100
86) Indeno(1,2,3-cd)pyrene	16.94	276	164877	45.568	nq	98
88) Benzo(b)fluoranthene	15.04	252	170024	41.825	nq	99
89) Benzo(k)fluoranthene	15.07	252	152391	39.574	nq	98
90) Benzo(a)pyrene	15.41	252	153336	42.985	nq	100
91) Dibenzo(a,h)anthracene	16.94	278	142007	49.972	nq	96
92) Benzo(g,h,i)perylene	17.37	276	134568	47.521	nq	96

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

