

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101017\
 Data File : BF099314.D
 Acq On : 10 Oct 2017 21:51
 Operator : SJ/JU
 Sample : I5745-03MSD
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HR-06-100617-AMSD

Manual Integrations
 APPROVED

Sohil
 10/11/2017 7:24:03 PM

Quant Time: Oct 11 01:41:12 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 05 17:20:48 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.87	152	89961	20.00	ng	0.00
21) Naphthalene-d8	8.15	136	358288	20.00	ng	0.00
38) Acenaphthene-d10	9.90	164	148280	20.00	ng	0.00
63) Phenanthrene-d10	11.39	188	218788	20.00	ng	0.00
75) Chrysene-d12	14.03	240	171723	20.00	ng	0.00
86) Perylene-d12	15.50	264	152319	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.51	112	726739	128.60	ng	0.03
7) Phenol-d6	6.50	99	803581	115.27	ng	0.00
23) Nitrobenzene-d5	7.43	82	547817	98.05	ng	0.00
41) 2,4,6-Tribromophenol	10.69	330	160217	132.05	ng	0.00
44) 2-Fluorobiphenyl	9.22	172	1013643	114.12	ng	0.00
78) Terphenyl-d14	12.97	244	684785	90.69	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.82	88	101189	37.484	ng	# 80
3) Pyridine	3.56	79	192595	26.373	ng	95
4) n-Nitrosodimethylamine	3.49	42	110854	35.797	ng	# 79
6) Aniline	6.53	93	140339	14.916	ng	98
8) 2-Chlorophenol	6.66	128	278899	43.580	ng	94
9) Benzaldehyde	6.42	77	95327	23.573	ng	91
10) Phenol	6.52	94	296788	38.066	ng	92
11) bis(2-Chloroethyl)ether	6.60	93	261255	43.103	ng	98
12) 1,3-Dichlorobenzene	6.81	146	298327	44.511	ng	96
13) 1,4-Dichlorobenzene	6.89	146	299117	43.864	ng	98
14) 1,2-Dichlorobenzene	7.04	146	279044	44.286	ng	97
15) Benzyl Alcohol	7.01	79	214264	41.896	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.14	45	365999	38.757	ng	94
17) 2-Methylphenol	7.12	107	226899	43.331	ng	96
18) Hexachloroethane	7.37	117	94847	38.696	ng	98
19) n-Nitroso-di-n-propylamine	7.28	70	177105	39.791	ng	96
20) 3+4-Methylphenols	7.27	107	269250	40.645	ng	# 70
22) Acetophenone	7.27	105	376738	43.400	ng	96
24) Nitrobenzene	7.45	77	274773	43.356	ng	96
25) Isophorone	7.69	82	498402	43.148	ng	98
26) 2-Nitrophenol	7.77	139	148984	48.885	ng	# 85
27) 2,4-Dimethylphenol	7.80	122	239788	41.663	ng	99
28) bis(2-Chloroethoxy)methane	7.89	93	326494	45.357	ng	99
29) 2,4-Dichlorophenol	8.01	162	229744	46.493	ng	96
30) 1,2,4-Trichlorobenzene	8.09	180	238134	48.183	ng	99
31) Naphthalene	8.17	128	797282	47.380	ng	99
32) Benzoic acid	7.92	122	94066	19.897	ng	95
33) 4-Chloroaniline	8.22	127	51642	7.349	ng	96
34) Hexachlorobutadiene	8.28	225	118249	46.452	ng	98
35) Caprolactam	8.60	113	59356m	37.446	ng	
36) 4-Chloro-3-methylphenol	8.70	107	228059	41.061	ng	95
37) 2-Methylnaphthalene	8.86	142	534560	48.866	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.03	216	215965	48.837	ng	98
40) Hexachlorocyclopentadiene	9.01	237	30704	15.193	ng	96

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101017\
 Data File : BF099314.D
 Acq On : 10 Oct 2017 21:51
 Operator : SJ/JU
 Sample : I5745-03MSD
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HR-06-100617-AMSD

Manual Integrations
 APPROVED

Sohil
 10/11/2017 7:24:03 PM

Quant Time: Oct 11 01:41:12 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 05 17:20:48 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.14	196	144998	46.284	nq	100
43) 2,4,5-Trichlorophenol	9.19	196	151663	48.496	nq	94
45) 1,1'-Biphenyl	9.32	154	609078	47.794	nq	99
46) 2-Chloronaphthalene	9.35	162	474158	49.555	nq	97
47) 2-Nitroaniline	9.45	65	128133	43.734	nq	87
48) Acenaphthylene	9.77	152	685867	46.825	nq	99
49) Dimethylphthalate	9.62	163	548484	49.010	nq	100
50) 2,6-Dinitrotoluene	9.69	165	118565	48.663	nq #	79
51) Acenaphthene	9.94	154	401606	43.284	nq	99
52) 3-Nitroaniline	9.86	138	61055	21.492	nq	94
53) 2,4-Dinitrophenol	9.97	184	22556	22.314	nq	93
54) Dibenzofuran	10.11	168	595453	48.200	nq	99
55) 4-Nitrophenol	10.03	139	147577	61.435	nq	85
56) 2,4-Dinitrotoluene	10.10	165	140385	44.397	nq #	87
57) Fluorene	10.45	166	451148	45.435	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.23	232	101734	47.092	nq	95
59) Diethylphthalate	10.32	149	501308	45.842	nq	97
60) 4-Chlorophenyl-phenylether	10.44	204	209349	46.935	nq	97
61) 4-Nitroaniline	10.47	138	109757	40.046	nq	87
62) Azobenzene	10.60	77	427114	42.478	nq	96
64) 4,6-Dinitro-2-methylphenol	10.50	198	19793	14.869	nq	84
65) n-Nitrosodiphenylamine	10.56	169	405462	53.495	nq	98
66) 4-Bromophenyl-phenylether	10.93	248	114632	53.527	nq	99
67) Hexachlorobenzene	11.00	284	111170	48.603	nq	97
68) Atrazine	11.09	200	117116	51.357	nq	98
69) Pentachlorophenol	11.20	266	103677	72.831	nq	99
70) Phenanthrene	11.42	178	575965	49.181	nq	98
71) Anthracene	11.47	178	587177	49.002	nq	99
72) Carbazole	11.62	167	538472	45.889	nq	99
73) Di-n-butylphthalate	11.94	149	663075	46.946	nq	99
74) Fluoranthene	12.60	202	562626	47.444	nq	98
76) Benzidine	12.72	184	194831	30.675	nq	98
77) Pyrene	12.83	202	584816	44.661	nq	98
79) Butylbenzylphthalate	13.44	149	257026	41.765	nq	95
80) Benzo(a)anthracene	14.02	228	507061	48.764	nq	99
81) 3,3'-Dichlorobenzidine	13.98	252	151795	39.822	nq	97
82) Chrysene	14.06	228	477133	48.197	nq	100
83) Bis(2-ethylhexyl)phthalate	13.99	149	348853	41.168	nq #	99
84) Di-n-octyl phthalate	14.60	149	630970	46.243	nq #	95
85) Indeno(1,2,3-cd)pyrene	17.00	276	386583	48.817	nq	97
87) Benzo(b)fluoranthene	15.07	252	454874	48.571	nq	97
88) Benzo(k)fluoranthene	15.10	252	459958	49.725	nq	99
89) Benzo(a)pyrene	15.44	252	426348	49.595	nq #	97
90) Dibenzo(a,h)anthracene	17.01	278	324626	47.185	nq	99
91) Benzo(g,h,i)perylene	17.45	276	312765	45.991	nq	97

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

