

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060117\
 Data File : BF095594.D
 Acq On : 1 Jun 2017 14:52
 Operator : SJ/MA
 Sample : PB99358BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB99358BS

Manual Integrations
 APPROVED

mohammad
 6/2/2017 2:32:20 PM

Quant Time: Jun 01 18:53:31 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 01 18:51:16 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.53	152	86621	20.00	ng	0.00
21) Naphthalene-d8	9.57	136	380073	20.00	ng	0.00
38) Acenaphthene-d10	12.40	164	170554	20.00	ng	0.00
63) Phenanthrene-d10	14.80	188	331158	20.00	ng	0.00
75) Chrysene-d12	18.50	240	249520	20.00	ng	0.00
86) Perylene-d12	20.17	264	167489	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.47	112	797373	153.47	ng	0.02
7) Phenol-d6	7.10	99	866231	138.96	ng	0.00
23) Nitrobenzene-d5	8.45	82	548522	91.01	ng	-0.01
41) 2,4,6-Tribromophenol	13.70	330	191304	136.96	ng	0.00
44) 2-Fluorobiphenyl	11.35	172	1003551	84.69	ng	0.00
78) Terphenyl-d14	17.15	244	946525	83.55	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.94	88	99362	38.995	ng	95
3) Pyridine	2.53	79	241937	40.968	ng	98
4) n-Nitrosodimethylamine	2.49	42	108955	44.262	ng	80
6) Aniline	7.04	93	181832	23.105	ng	98
8) 2-Chlorophenol	7.23	128	263753	44.601	ng	95
9) Benzaldehyde	6.86	77	36569	9.607	ng	91
10) Phenol	7.12	94	328137	49.952	ng	90
11) bis(2-Chloroethyl)ether	7.17	93	222939	41.109	ng	97
12) 1,3-Dichlorobenzene	7.43	146	275778	41.111	ng	97
13) 1,4-Dichlorobenzene	7.56	146	280368	41.213	ng	97
14) 1,2-Dichlorobenzene	7.80	146	272735	42.951	ng	96
15) Benzyl Alcohol	7.83	79	219325	54.701	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.04	45	314495	40.973	ng	96
17) 2-Methylphenol	8.06	107	233183	48.183	ng	94
18) Hexachloroethane	8.32	117	101224	43.924	ng	87
19) n-Nitroso-di-n-propylamine	8.26	70	196589	46.628	ng	94
20) 3+4-Methylphenols	8.33	107	317516	48.283	ng	# 60
22) Acetophenone	8.22	105	393461	43.147	ng	# 83
24) Nitrobenzene	8.48	77	276751	43.806	ng	91
25) Isophorone	8.88	82	489292	45.463	ng	95
26) 2-Nitrophenol	8.99	139	155471	45.606	ng	98
27) 2,4-Dimethylphenol	9.14	122	239902	43.527	ng	98
28) bis(2-Chloroethoxy)methane	9.26	93	318210	42.740	ng	98
29) 2,4-Dichlorophenol	9.42	162	239005	45.926	ng	97
30) 1,2,4-Trichlorobenzene	9.50	180	228537	40.990	ng	96
31) Naphthalene	9.60	128	777897	40.723	ng	100
32) Benzoic acid	9.47	122	94993	28.553	ng	91
33) 4-Chloroaniline	9.76	127	132970	18.679	ng	# 90
34) Hexachlorobutadiene	9.82	225	105985	36.026	ng	99
35) Caprolactam	10.39	113	74024m	47.369	ng	
36) 4-Chloro-3-methylphenol	10.62	107	235548	42.334	ng	91
37) 2-Methylnaphthalene	10.73	142	512802	40.903	ng	97
39) 1,2,4,5-Tetrachlorobenzene	11.01	216	200000	38.132	ng	100
40) Hexachlorocyclopentadiene	10.98	237	100174	48.048	ng	98

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF060117\
 Data File : BF095594.D
 Acq On : 1 Jun 2017 14:52
 Operator : SJ/MA
 Sample : PB99358BS
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleID :
 PB99358BS

Manual Integrations
 APPROVED

mohammad
 6/2/2017 2:32:20 PM

Quant Time: Jun 01 18:53:31 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 01 18:51:16 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	11.24	196	147391	42.089	ng	98
43) 2,4,5-Trichlorophenol	11.33	196	140712	37.385	ng	94
45) 1,1'-Biphenyl	11.50	154	596716	41.555	ng	97
46) 2-Chloronaphthalene	11.51	162	467237	40.297	ng	94
47) 2-Nitroaniline	11.73	65	142823	45.004	ng	# 79
48) Acenaphthylene	12.17	152	819800	45.141	ng	98
49) Dimethylphthalate	12.05	163	581188	41.509	ng	99
50) 2,6-Dinitrotoluene	12.14	165	132572	46.411	ng	# 87
51) Acenaphthene	12.45	154	448995	41.392	ng	98
52) 3-Nitroaniline	12.42	138	85892	28.015	ng	90
53) 2,4-Dinitrophenol	12.63	184	73210	49.146	ng	# 65
54) Dibenzofuran	12.74	168	700799	46.686	ng	97
55) 4-Nitrophenol	12.82	139	145105	78.800	ng	# 74
56) 2,4-Dinitrotoluene	12.82	165	177824	48.447	ng	# 87
57) Fluorene	13.29	166	589395	45.156	ng	99
58) 2,3,4,6-Tetrachlorophenol	12.99	232	119565	50.475	ng	# 92
59) Diethylphthalate	13.20	149	564780	42.084	ng	99
60) 4-Chlorophenyl-phenylether	13.32	204	255696	44.574	ng	92
61) 4-Nitroaniline	13.42	138	132641	47.127	ng	97
62) Azobenzene	13.57	77	517054	47.947	ng	95
64) 4,6-Dinitro-2-methylphenol	13.49	198	73728	33.211	ng	# 68
65) n-Nitrosodiphenylamine	13.53	169	513507	42.725	ng	98
66) 4-Bromophenyl-phenylether	14.10	248	141351	40.402	ng	95
67) Hexachlorobenzene	14.19	284	141069	39.344	ng	# 89
68) Atrazine	14.46	200	153594	45.261	ng	98
69) Pentachlorophenol	14.55	266	83773	53.750	ng	96
70) Phenanthrene	14.84	178	827031	43.465	ng	98
71) Anthracene	14.92	178	825556	42.476	ng	98
72) Carbazole	15.22	167	782662	44.258	ng	97
73) Di-n-butylphthalate	15.79	149	884189	40.093	ng	99
74) Fluoranthene	16.60	202	832393	42.647	ng	99
76) Benzidine	16.83	184	164970	27.679	ng	# 92
77) Pyrene	16.90	202	836506	44.826	ng	99
79) Butylbenzylphthalate	17.81	149	384220	44.265	ng	# 86
80) Benzo(a)anthracene	18.49	228	621162	42.774	ng	99
81) 3,3'-Dichlorobenzidine	18.47	252	133762	26.669	ng	# 96
82) Chrysene	18.53	228	592795	42.217	ng	99
83) Bis(2-ethylhexyl)phthalate	18.56	149	520812	42.112	ng	98
84) Di-n-octyl phthalate	19.33	149	787821	38.855	ng	96
85) Indeno(1,2,3-cd)pyrene	21.40	276	413913	36.298	ng	100
87) Benzo(b)fluoranthene	19.76	252	477106	46.100	ng	98
88) Benzo(k)fluoranthene	19.79	252	441019	44.177	ng	96
89) Benzo(a)pyrene	20.11	252	415927	43.553	ng	98
90) Dibenzo(a,h)anthracene	21.40	278	341935	43.354	ng	97
91) Benzo(g,h,i)perylene	21.76	276	366104	47.301	ng	98

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

