

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF091217\
 Data File : BF098451.D
 Acq On : 12 Sep 2017 15:40
 Operator : SJ/JU
 Sample : I5156-01MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-02-011-AMS

Manual Integrations
 APPROVED

Sohil
 9/13/2017 4:59:58 PM

Quant Time: Sep 13 01:15:09 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF091117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 12 02:14:50 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.66	152	222845	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	940781	20.00	ng	0.00
38) Acenaphthene-d10	9.70	164	419236	20.00	ng	0.00
63) Phenanthrene-d10	11.18	188	709094	20.00	ng	0.00
75) Chrysene-d12	13.81	240	402857	20.00	ng	0.00
86) Perylene-d12	15.21	264	392465	20.00	ng	0.00

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System Monitoring Compounds

5) 2-Fluorophenol	5.28	112	1469893	97.52	ng	0.00
7) Phenol-d6	6.32	99	1816272	94.91	ng	0.00
23) Nitrobenzene-d5	7.23	82	1132780	67.13	ng	0.00
41) 2,4,6-Tribromophenol	10.49	330	305031	109.02	ng	0.00
44) 2-Fluorobiphenyl	9.02	172	1757054	71.04	ng	-0.01
78) Terphenyl-d14	12.76	244	1188414	63.67	ng	-0.01

9/13/2017 4:59:58 PM

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.34	88	207624	26.689	ng	89
3) Pyridine	3.03	79	592335	28.672	ng	# 83
4) n-Nitrosodimethylamine	2.99	42	295342	29.160	ng	100
6) Aniline	6.33	93	521581	21.726	ng	# 1
8) 2-Chlorophenol	6.45	128	546977	33.003	ng	94
9) Benzaldehyde	6.21	77	410803	32.839	ng	91
10) Phenol	6.33	94	760640	34.221	ng	89
11) bis(2-Chloroethyl)ether	6.40	93	538214	32.115	ng	90
12) 1,3-Dichlorobenzene	6.60	146	550579	33.005	ng	98
13) 1,4-Dichlorobenzene	6.68	146	549068	32.133	ng	94
14) 1,2-Dichlorobenzene	6.83	146	506968	32.566	ng	98
15) Benzyl Alcohol	6.82	79	451844	35.478	ng	96
16) 2,2'-oxybis(1-Chloropropan	6.95	45	877305	31.816	ng	74
17) 2-Methylphenol	6.94	107	459354	33.989	ng	# 89
18) Hexachloroethane	7.17	117	208424	31.851	ng	# 81
19) n-Nitroso-di-n-propylamine	7.09	70	391254	32.504	ng	99
20) 3+4-Methylphenols	7.09	107	602190	35.309	ng	# 66
22) Acetophenone	7.08	105	772365	31.765	ng	# 92
24) Nitrobenzene	7.25	77	608428	31.653	ng	98
25) Isophorone	7.49	82	1083041	31.721	ng	100
26) 2-Nitrophenol	7.57	139	286667	36.937	ng	# 83
27) 2,4-Dimethylphenol	7.62	122	489792	33.106	ng	96
28) bis(2-Chloroethoxy)methane	7.70	93	696576	33.435	ng	98
29) 2,4-Dichlorophenol	7.82	162	436051	35.438	ng	96
30) 1,2,4-Trichlorobenzene	7.89	180	456368	34.096	ng	99
31) Naphthalene	7.97	128	1549377	33.314	ng	99
32) Benzoic acid	7.72	122	18625	9.535	ng	93
33) 4-Chloroaniline	8.02	127	327550	17.329	ng	99
34) Hexachlorobutadiene	8.09	225	230898	33.603	ng	99
35) Caprolactam	8.40	113	155444m	36.635	ng	
36) 4-Chloro-3-methylphenol	8.52	107	497531	32.604	ng	# 82
37) 2-Methylnaphthalene	8.66	142	1016203	36.072	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.83	216	416425	31.888	ng	99
40) Hexachlorocyclopentadiene	8.81	237	128782	39.368	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.95	196	276597	36.060	nq	92
43) 2,4,5-Trichlorophenol	8.99	196	290624	36.249	nq	93
45) 1,1'-Biphenyl	9.12	154	1190067	31.715	nq	97
46) 2-Chloronaphthalene	9.15	162	899751	33.574	nq	94
47) 2-Nitroaniline	9.25	65	352394	34.260	nq	93
48) Acenaphthylene	9.56	152	1321649	33.169	nq	99
49) Dimethylphthalate	9.43	163	1259515	38.811	nq	99
50) 2,6-Dinitrotoluene	9.50	165	236718	35.511	nq	# 85
51) Acenaphthene	9.73	154	822286	32.465	nq	98
52) 3-Nitroaniline	9.66	138	193621	24.062	nq	98
53) 2,4-Dinitrophenol	9.78	184	60451	25.665	nq	# 76
54) Dibenzofuran	9.90	168	1155764	34.637	nq	98
55) 4-Nitrophenol	9.85	139	304524	65.870	nq	# 46
56) 2,4-Dinitrotoluene	9.90	165	285577	35.238	nq	# 58
57) Fluorene	10.25	166	893978	32.369	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.03	232	212639	39.464	nq	94
59) Diethylphthalate	10.13	149	1028090	33.311	nq	98
60) 4-Chlorophenyl-phenylether	10.24	204	428882	33.620	nq	# 87
61) 4-Nitroaniline	10.28	138	242209	29.805	nq	88
62) Azobenzene	10.40	77	1074367	32.298	nq	96
64) 4,6-Dinitro-2-methylphenol	10.31	198	102558	26.106	nq	89
65) n-Nitrosodiphenylamine	10.36	169	832338	36.150	nq	99
66) 4-Bromophenyl-phenylether	10.73	248	247796	36.820	nq	95
67) Hexachlorobenzene	10.79	284	232096	34.000	nq	# 86
68) Atrazine	10.89	200	260100	36.694	nq	98
69) Pentachlorophenol	10.99	266	178274	60.818	nq	95
70) Phenanthrene	11.20	178	1310793	34.870	nq	99
71) Anthracene	11.26	178	1321941	34.252	nq	99
72) Carbazole	11.42	167	1125435	31.289	nq	99
73) Di-n-butylphthalate	11.75	149	1485821	34.477	nq	100
74) Fluoranthene	12.39	202	1202466	31.954	nq	98
76) Benzidine	12.52	184	471741	26.459	nq	# 92
77) Pyrene	12.62	202	1196604	37.653	nq	99
79) Butylbenzylphthalate	13.23	149	500571	36.308	nq	# 76
80) Benzo(a)anthracene	13.80	228	851825	36.066	nq	97
81) 3,3'-Dichlorobenzidine	13.77	252	244939	26.685	nq	# 94
82) Chrysene	13.84	228	814041	34.096	nq	100
83) Bis(2-ethylhexyl)phthalate	13.79	149	629224	36.365	nq	100
84) Di-n-octyl phthalate	14.40	149	1078160	36.318	nq	97
85) Indeno(1,2,3-cd)pyrene	16.56	276	679418	40.562	nq	98
87) Benzo(b)fluoranthene	14.82	252	897110	36.354	nq	97
88) Benzo(k)fluoranthene	14.84	252	702006	30.472	nq	96
89) Benzo(a)pyrene	15.16	252	784154	35.668	nq	99
90) Dibenzo(a,h)anthracene	16.56	278	564060	35.270	nq	98
91) Benzo(g,h,i)perylene	16.96	276	547877	34.061	nq	100

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

