

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF091217\
 Data File : BF098452.D
 Acq On : 12 Sep 2017 16:07
 Operator : SJ/JU
 Sample : I5156-01MSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 9/13/2017 5:00:01 PM

Quant Time: Sep 13 01:21:32 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	215104	20.00	ng	0.00
21) Naphthalene-d8	7.95	136	914049	20.00	ng	0.00
38) Acenaphthene-d10	9.70	164	399310	20.00	ng	0.00
63) Phenanthrene-d10	11.18	188	649078	20.00	ng	0.00
75) Chrysene-d12	13.81	240	385079	20.00	ng	0.00
86) Perylene-d12	15.21	264	385418	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.28	112	1531282	105.25	ng	0.00
7) Phenol-d6	6.32	99	1880628	101.81	ng	0.00
23) Nitrobenzene-d5	7.23	82	1193484	72.79	ng	0.00
41) 2,4,6-Tribromophenol	10.49	330	312993	117.44	ng	0.00
44) 2-Fluorobiphenyl	9.02	172	1811964	76.91	ng	-0.01
78) Terphenyl-d14	12.76	244	1216824	68.21	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.35	88	215150	28.652	ng	87
3) Pyridine	3.03	79	546594	27.410	ng	# 80
4) n-Nitrosodimethylamine	2.99	42	308705	31.576	ng	99
6) Aniline	6.33	93	538699	23.246	ng	# 6
8) 2-Chlorophenol	6.45	128	563389	35.217	ng	93
9) Benzaldehyde	6.21	77	418017	34.618	ng	92
10) Phenol	6.33	94	756131	35.242	ng	87
11) bis(2-Chloroethyl)ether	6.40	93	558591	34.531	ng	91
12) 1,3-Dichlorobenzene	6.60	146	556469	34.559	ng	97
13) 1,4-Dichlorobenzene	6.68	146	560410	33.977	ng	95
14) 1,2-Dichlorobenzene	6.83	146	521180	34.684	ng	98
15) Benzyl Alcohol	6.82	79	461151	37.512	ng	95
16) 2,2'-oxybis(1-Chloropropan	6.95	45	903198	33.934	ng	72
17) 2-Methylphenol	6.94	107	474697	36.389	ng	# 90
18) Hexachloroethane	7.17	117	210763	33.367	ng	# 82
19) n-Nitroso-di-n-propylamine	7.09	70	406831	35.014	ng	97
20) 3+4-Methylphenols	7.09	107	618652	37.579	ng	# 67
22) Acetophenone	7.08	105	806475	34.138	ng	# 91
24) Nitrobenzene	7.25	77	624350	33.431	ng	98
25) Isophorone	7.49	82	1112467	33.536	ng	100
26) 2-Nitrophenol	7.57	139	294352	39.036	ng	# 83
27) 2,4-Dimethylphenol	7.62	122	505824	35.189	ng	97
28) bis(2-Chloroethoxy)methane	7.70	93	710811	35.116	ng	99
29) 2,4-Dichlorophenol	7.82	162	441838	36.959	ng	96
30) 1,2,4-Trichlorobenzene	7.89	180	468385	36.017	ng	99
31) Naphthalene	7.97	128	1580802	34.984	ng	100
32) Benzoic acid	7.71	122	8584m	8.712	ng	
33) 4-Chloroaniline	8.02	127	369977	20.146	ng	99
34) Hexachlorobutadiene	8.09	225	232236	34.786	ng	98
35) Caprolactam	8.40	113	137732m	33.410	ng	
36) 4-Chloro-3-methylphenol	8.52	107	505912	34.123	ng	# 83
37) 2-Methylnaphthalene	8.66	142	1027460	37.538	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.83	216	419628	33.737	ng	99
40) Hexachlorocyclopentadiene	8.81	237	118190	38.291	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.95	196	278439	38.111	nq	92
43) 2,4,5-Trichlorophenol	8.99	196	289490	37.909	nq	93
45) 1,1'-Biphenyl	9.12	154	1213936	33.965	nq	97
46) 2-Chloronaphthalene	9.15	162	914941	35.845	nq #	92
47) 2-Nitroaniline	9.25	65	354481	36.183	nq	95
48) Acenaphthylene	9.56	152	1331348	35.080	nq	99
49) Dimethylphthalate	9.43	163	1206153	39.022	nq	99
50) 2,6-Dinitrotoluene	9.50	165	240151	37.824	nq #	85
51) Acenaphthene	9.73	154	841834	34.895	nq	99
52) 3-Nitroaniline	9.66	138	202743	26.453	nq	100
53) 2,4-Dinitrophenol	9.78	184	28384	16.974	nq #	79
54) Dibenzofuran	9.90	168	1146003	36.058	nq	99
55) 4-Nitrophenol	9.85	139	301846	68.548	nq #	44
56) 2,4-Dinitrotoluene	9.90	165	281034	36.408	nq #	57
57) Fluorene	10.25	166	898900	34.171	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.03	232	200568	39.082	nq	99
59) Diethylphthalate	10.13	149	1031269	35.082	nq	98
60) 4-Chlorophenyl-phenylether	10.24	204	432122	35.564	nq #	87
61) 4-Nitroaniline	10.28	138	243076	31.404	nq	91
62) Azobenzene	10.40	77	1083073	34.185	nq	97
64) 4,6-Dinitro-2-methylphenol	10.31	198	78384	22.617	nq	86
65) n-Nitrosodiphenylamine	10.36	169	830212	39.392	nq	98
66) 4-Bromophenyl-phenylether	10.73	248	251370	40.805	nq	94
67) Hexachlorobenzene	10.79	284	231858	37.105	nq #	86
68) Atrazine	10.89	200	255987	39.453	nq	98
69) Pentachlorophenol	10.99	266	157121	58.815	nq	96
70) Phenanthrene	11.20	178	1283815	37.310	nq	99
71) Anthracene	11.26	178	1293046	36.601	nq	99
72) Carbazole	11.42	167	1102430	33.484	nq	99
73) Di-n-butylphthalate	11.74	149	1456072	36.911	nq	99
74) Fluoranthene	12.39	202	1113512	32.326	nq	98
76) Benzidine	12.52	184	519986	30.511	nq #	93
77) Pyrene	12.62	202	1111032	36.574	nq	99
79) Butylbenzylphthalate	13.23	149	502889	38.160	nq #	75
80) Benzo(a)anthracene	13.80	228	802435	35.543	nq	98
81) 3,3'-Dichlorobenzidine	13.76	252	265469	30.257	nq	98
82) Chrysene	13.84	228	816676	35.786	nq	99
83) Bis(2-ethylhexyl)phthalate	13.79	149	633686	38.314	nq	99
84) Di-n-octyl phthalate	14.40	149	1100339	38.776	nq	97
85) Indeno(1,2,3-cd)pyrene	16.56	276	669305	41.803	nq	98
87) Benzo(b)fluoranthene	14.82	252	799433	32.988	nq	97
88) Benzo(k)fluoranthene	14.84	252	794643	35.124	nq	96
89) Benzo(a)pyrene	15.16	252	770979	35.710	nq	99
90) Dibenzo(a,h)anthracene	16.56	278	567631	36.142	nq	96
91) Benzo(g,h,i)perylene	16.96	276	540760	34.233	nq	99

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

