

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF122617\
 Data File : BF101744.D
 Acq On : 27 Dec 2017 5:07
 Operator : SJ/JU
 Sample : I7043-01MSD
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EPE-106-3DS(90-92)MSD

Manual Integrations
 APPROVED

Sohil
 12/27/2017 4:15:25 PM

Quant Time: Dec 27 07:20:42 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 26 10:48:01 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	135013	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	487927	20.00	ng	0.00
38) Acenaphthene-d10	9.83	164	185041	20.00	ng	0.00
63) Phenanthrene-d10	11.32	188	298800	20.00	ng	0.00
75) Chrysene-d12	13.97	240	287240	20.00	ng	0.00
86) Perylene-d12	15.44	264	262873	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.42	112	766183	84.53	ng	0.01
7) Phenol-d6	6.44	99	908020	80.50	ng	0.00
23) Nitrobenzene-d5	7.36	82	583036	66.52	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	222011	124.51	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	1008260	84.52	ng	0.00
78) Terphenyl-d14	12.90	244	1113395	93.45	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.57	88	105244	22.598	ng	97
3) Pyridine	3.35	79	237058	18.320	ng	95
4) n-Nitrosodimethylamine	3.29	42	123384	20.709	ng	97
6) Aniline	6.47	93	83705	5.340	ng	# 53
8) 2-Chlorophenol	6.58	128	218655	22.933	ng	98
9) Benzaldehyde	6.36	77	77324	9.282	ng	98
10) Phenol	6.45	94	306787	23.862	ng	99
11) bis(2-Chloroethyl)ether	6.53	93	240036	23.801	ng	97
12) 1,3-Dichlorobenzene	6.73	146	252633	23.477	ng	99
13) 1,4-Dichlorobenzene	6.80	146	261623	23.808	ng	99
14) 1,2-Dichlorobenzene	6.96	146	234499	23.708	ng	97
15) Benzyl Alcohol	6.95	79	200985	25.886	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.06	45	409090	26.505	ng	# 44
17) 2-Methylphenol	7.06	107	199609	22.759	ng	# 87
18) Hexachloroethane	7.29	117	70284	18.396	ng	94
19) n-Nitroso-di-n-propylamine	7.20	70	205536	27.745	ng	96
20) 3+4-Methylphenols	7.22	107	297408	32.000	ng	99
22) Acetophenone	7.20	105	388491	34.894	ng	# 97
24) Nitrobenzene	7.38	77	267900	27.616	ng	99
25) Isophorone	7.62	82	550287	31.295	ng	99
26) 2-Nitrophenol	7.70	139	114463	27.464	ng	94
27) 2,4-Dimethylphenol	7.73	122	236613	33.418	ng	98
28) bis(2-Chloroethoxy)methane	7.82	93	340594	30.493	ng	98
29) 2,4-Dichlorophenol	7.96	162	213096	30.464	ng	98
30) 1,2,4-Trichlorobenzene	8.02	180	194679	26.372	ng	97
31) Naphthalene	8.09	128	688987	28.049	ng	100
32) Benzoic acid	7.86	122	37212m	14.467	ng	
33) 4-Chloroaniline	8.17	127	57440	5.380	ng	# 93
34) Hexachlorobutadiene	8.20	225	110739	25.937	ng	99
35) Caprolactam	8.53	113	90795	37.381	ng	93
36) 4-Chloro-3-methylphenol	8.65	107	231731	30.140	ng	99
37) 2-Methylnaphthalene	8.79	142	470770	29.416	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.95	216	221712	35.488	ng	# 97
42) 2,4,6-Trichlorophenol	9.07	196	137810	36.640	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	9.13	196	131530	31.939	ng	95
45) 1,1'-Biphenyl	9.25	154	577251	35.176	ng	99
46) 2-Chloronaphthalene	9.27	162	420998	33.528	ng	97
47) 2-Nitroaniline	9.39	65	163690	37.737	ng	91
48) Acenaphthylene	9.69	152	611200	30.818	ng	100
49) Dimethylphthalate	9.55	163	648960	43.601	ng	98
50) 2,6-Dinitrotoluene	9.63	165	107686	35.598	ng	91
51) Acenaphthene	9.86	154	376157	31.569	ng	100
52) 3-Nitroaniline	9.80	138	81387	21.579	ng	99
54) Dibenzofuran	10.03	168	556867	36.775	ng	97
55) 4-Nitrophenol	9.99	139	111764	67.960	ng	94
56) 2,4-Dinitrotoluene	10.05	165	141665	41.565	ng #	96
57) Fluorene	10.38	166	452789	35.347	ng	98
58) 2,3,4,6-Tetrachlorophenol	10.17	232	115419	39.009	ng #	86
59) Diethylphthalate	10.24	149	570958	38.737	ng	96
60) 4-Chlorophenyl-phenylether	10.36	204	221092	36.013	ng	94
61) 4-Nitroaniline	10.42	138	130296	34.370	ng	94
62) Azobenzene	10.53	77	484741	31.747	ng	96
65) n-Nitrosodiphenylamine	10.49	169	442875	40.141	ng	96
66) 4-Bromophenyl-phenylether	10.86	248	137367	40.915	ng #	83
67) Hexachlorobenzene	10.93	284	142729	40.167	ng #	87
68) Atrazine	11.02	200	104677	31.819	ng	95
69) Pentachlorophenol	11.13	266	117620	84.700	ng	100
70) Phenanthrene	11.34	178	678817	40.851	ng	99
71) Anthracene	11.40	178	713207	42.019	ng	98
72) Carbazole	11.56	167	701632	44.151	ng	100
73) Di-n-butylphthalate	11.87	149	873195	45.271	ng	99
74) Fluoranthene	12.54	202	795860	45.630	ng	97
76) Benzidine	12.66	184	232684	19.180	ng	97
77) Pyrene	12.77	202	838736	38.907	ng	99
79) Butylbenzylphthalate	13.37	149	392863	40.276	ng	97
80) Benzo(a)anthracene	13.96	228	816673	43.058	ng	99
81) 3,3'-Dichlorobenzidine	13.92	252	213902	34.587	ng	98
82) Chrysene	14.00	228	795023	44.394	ng	100
83) Bis(2-ethylhexyl)phthalate	13.92	149	527418	42.689	ng	98
84) Di-n-octyl phthalate	14.53	149	1023654	46.459	ng	99
85) Indeno(1,2,3-cd)pyrene	16.92	276	614229	38.516	ng	97
87) Benzo(b)fluoranthene	15.02	252	749392	46.771	ng	99
88) Benzo(k)fluoranthene	15.04	252	688978	44.226	ng	98
89) Benzo(a)pyrene	15.38	252	669087	45.298	ng	100
90) Dibenzo(a,h)anthracene	16.92	278	524329	41.345	ng	97
91) Benzo(g,h,i)perylene	17.37	276	514410	40.964	ng	95

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

