

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011316\
 Data File : BF084475.D
 Acq On : 14 Jan 2016 7:49
 Operator : SJ/UM
 Sample : H1079-02MSD
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEC-COMP1MSD

Quant Time: Jan 14 16:30:36 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 14 14:15:52 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	265783	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	1087335	20.00	ng	0.00
38) Acenaphthene-d10	9.87	164	475127	20.00	ng	0.00
63) Phenanthrene-d10	11.36	188	787768	20.00	ng	0.00
75) Chrysene-d12	14.00	240	490551	20.00	ng	0.00
86) Perylene-d12	15.41	264	498440	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.45	112	1998506	121.62	ng	0.01
7) Phenol-d6	6.49	99	2674210	129.58	ng	0.01
23) Nitrobenzene-d5	7.40	82	1621286	82.99	ng	0.00
41) 2,4,6-Tribromophenol	10.67	330	608292	126.81	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	2315509	85.79	ng	0.00
78) Terphenyl-d14	12.95	244	1932491	87.94	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.50	88	318634	40.93	ng	# 77
3) Pyridine	3.21	79	930447	42.87	ng	# 77
4) n-Nitrosodimethylamine	3.16	42	455331	40.87	ng	87
6) Aniline	6.50	93	597750	19.80	ng	# 1
8) 2-Chlorophenol	6.62	128	861172	46.19	ng	97
9) Benzaldehyde	6.37	77	184097	13.70	ng	96
10) Phenol	6.50	94	1182987	50.42	ng	95
11) bis(2-Chloroethyl)ether	6.58	93	883890	48.63	ng	91
12) 1,3-Dichlorobenzene	6.77	146	827287m	43.02	ng	
13) 1,4-Dichlorobenzene	6.85	146	879253	45.59	ng	94
14) 1,2-Dichlorobenzene	7.00	146	735945	39.85	ng	95
15) Benzyl Alcohol	6.98	79	679389	39.13	ng	94
16) 2,2'-oxybis(1-Chloropropan	7.12	45	1301820	39.33	ng	88
17) 2-Methylphenol	7.10	107	727015	46.53	ng	95
18) Hexachloroethane	7.34	117	315699	40.94	ng	96
19) n-Nitroso-di-n-propylamine	7.25	70	553832	39.64	ng	# 75
20) 3+4-Methylphenols	7.25	107	796121	43.35	ng	# 60
22) Acetophenone	7.25	105	1141630	43.66	ng	# 93
24) Nitrobenzene	7.42	77	978340	49.37	ng	95
25) Isophorone	7.66	82	1644480	44.01	ng	97
26) 2-Nitrophenol	7.73	139	448078	49.69	ng	# 80
27) 2,4-Dimethylphenol	7.78	122	793584	43.47	ng	94
28) bis(2-Chloroethoxy)methane	7.87	93	1043983	45.74	ng	99
29) 2,4-Dichlorophenol	7.98	162	690863	45.43	ng	99
30) 1,2,4-Trichlorobenzene	8.06	180	718392	45.76	ng	# 96
31) Naphthalene	8.14	128	2458992	49.85	ng	98
32) Benzoic acid	7.86	122	56361	10.42	ng	94
33) 4-Chloroaniline	8.19	127	331958	14.68	ng	99
34) Hexachlorobutadiene	8.26	225	365599	40.52	ng	98
35) Caprolactam	8.57	113	22405	4.37	ng	# 57
36) 4-Chloro-3-methylphenol	8.68	107	711874	41.80	ng	96
37) 2-Methylnaphthalene	8.84	142	1966688	55.54	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.00	216	626952	45.90	ng	98
40) Hexachlorocyclopentadiene	8.99	237	484428	58.75	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.12	196	464200	45.42	nq	96
43) 2,4,5-Trichlorophenol	9.16	196	466050	47.57	nq	96
45) 1,1'-Biphenyl	9.30	154	1703013	45.10	nq	99
46) 2-Chloronaphthalene	9.32	162	1263814	42.61	nq	95
47) 2-Nitroaniline	9.42	65	525458	53.67	nq #	81
48) Acenaphthylene	9.73	152	1968740	44.93	nq	97
49) Dimethylphthalate	9.61	163	1994415	58.72	nq #	90
50) 2,6-Dinitrotoluene	9.66	165	339216	45.53	nq #	58
51) Acenaphthene	9.90	154	1327125	45.15	nq	98
52) 3-Nitroaniline	9.84	138	329333	35.68	nq #	82
53) 2,4-Dinitrophenol	9.94	184	146617	47.70	nq	86
54) Dibenzofuran	10.08	168	1748117	45.98	nq	91
55) 4-Nitrophenol	10.02	139	622978	92.97	nq	90
56) 2,4-Dinitrotoluene	10.06	165	417812	44.31	nq #	87
57) Fluorene	10.42	166	1214690	41.05	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.20	232	373111	44.61	nq	90
59) Diethylphthalate	10.30	149	1465761	43.77	nq	99
60) 4-Chlorophenyl-phenylether	10.42	204	636598	52.58	nq	94
61) 4-Nitroaniline	10.45	138	306004	37.19	nq #	74
62) Azobenzene	10.58	77	1689047	45.98	nq	95
64) 4,6-Dinitro-2-methylphenol	10.48	198	169618	35.35	nq	81
65) n-Nitrosodiphenylamine	10.53	169	1209091	48.00	nq	99
66) 4-Bromophenyl-phenylether	10.91	248	384337	45.33	nq	98
67) Hexachlorobenzene	10.97	284	372739	39.06	nq #	81
68) Atrazine	11.08	200	307632	37.32	nq	92
69) Pentachlorophenol	11.17	266	392798	79.38	nq	97
70) Phenanthrene	11.39	178	2013036	48.85	nq	97
71) Anthracene	11.44	178	1806343	44.72	nq	97
72) Carbazole	11.60	167	1867333	47.29	nq	98
73) Di-n-butylphthalate	11.93	149	2195096	55.37	nq #	97
74) Fluoranthene	12.57	202	1722262	40.01	nq	99
76) Benzidine	12.69	184	513802	29.21	nq	98
77) Pyrene	12.80	202	1731034	44.97	nq	98
79) Butylbenzylphthalate	13.43	149	768661	46.14	nq #	82
80) Benzo(a)anthracene	13.99	228	1250221	43.40	nq	99
81) 3,3'-Dichlorobenzidine	13.95	252	323573	32.19	nq #	98
82) Chrysene	14.02	228	1217876	44.00	nq	98
83) Bis(2-ethylhexyl)phthalate	13.99	149	1000637	47.55	nq #	94
84) Di-n-octyl phthalate	14.59	149	1589066	44.72	nq #	85
85) Indeno(1,2,3-cd)pyrene	16.80	276	1346333	61.36	nq	98
87) Benzo(b)fluoranthene	15.00	252	1492368m	45.77	nq	
88) Benzo(k)fluoranthene	15.04	252	955689m	35.84	nq	
89) Benzo(a)pyrene	15.36	252	1217309	44.02	nq #	96
90) Dibenzo(a,h)anthracene	16.82	278	1091797	45.88	nq	97
91) Benzo(g,h,i)perylene	17.22	276	1129124	45.82	nq	98

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

