

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052416\
 Data File : BF087348.D
 Acq On : 24 May 2016 20:53
 Operator : UM/SJ
 Sample : H3135-03MS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SEWER-JOBMS

Quant Time: May 25 15:26:51 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF052316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 25 00:58:25 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	109653	20.00	ng	-0.01
21) Naphthalene-d8	9.59	136	455137	20.00	ng	-0.01
38) Acenaphthene-d10	12.42	164	218373	20.00	ng	0.00
63) Phenanthrene-d10	14.82	188	432768	20.00	ng	0.00
75) Chrysene-d12	18.51	240	343512	20.00	ng	0.00
86) Perylene-d12	20.18	264	241151	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	996345	159.66	ng	0.03
7) Phenol-d6	7.11	99	1246001	147.77	ng	0.00
23) Nitrobenzene-d5	8.48	82	935104	103.55	ng	-0.01
41) 2,4,6-Tribromophenol	13.73	330	351605	167.55	ng	0.00
44) 2-Fluorobiphenyl	11.37	172	1605097	103.62	ng	-0.01
78) Terphenyl-d14	17.18	244	1585083	98.10	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.06	88	126804m	38.51	ng	
3) Pyridine	2.71	79	221460m	30.77	ng	
4) n-Nitrosodimethylamine	2.57	42	279095	50.32	ng	# 46
6) Aniline	7.07	93	136461	12.51	ng	95
8) 2-Chlorophenol	7.24	128	402646	51.74	ng	90
9) Benzaldehyde	6.91	77	40155	8.63	ng	95
10) Phenol	7.13	94	487201	52.91	ng	78
11) bis(2-Chloroethyl)ether	7.21	93	413875	55.53	ng	91
12) 1,3-Dichlorobenzene	7.46	146	415000	48.89	ng	97
13) 1,4-Dichlorobenzene	7.59	146	425385	48.82	ng	96
14) 1,2-Dichlorobenzene	7.82	146	407979	50.61	ng	98
15) Benzyl Alcohol	7.86	79	387450	63.02	ng	91
16) 2,2'-oxybis(1-Chloropropan	8.06	45	795758	47.46	ng	98
17) 2-Methylphenol	8.07	107	333626	53.49	ng	# 84
18) Hexachloroethane	8.33	117	158835	48.86	ng	96
19) n-Nitroso-di-n-propylamine	8.28	70	348705	55.45	ng	# 77
20) 3+4-Methylphenols	8.33	107	432120	57.31	ng	# 60
22) Acetophenone	8.24	105	604006	55.55	ng	# 97
24) Nitrobenzene	8.50	77	487380	51.13	ng	98
25) Isophorone	8.90	82	861219	53.17	ng	99
26) 2-Nitrophenol	9.00	139	213779	54.86	ng	# 70
27) 2,4-Dimethylphenol	9.15	122	370869	54.83	ng	89
28) bis(2-Chloroethoxy)methane	9.28	93	517575	56.73	ng	98
29) 2,4-Dichlorophenol	9.42	162	344314	56.48	ng	95
30) 1,2,4-Trichlorobenzene	9.51	180	359414	51.50	ng	96
31) Naphthalene	9.62	128	1317521	57.77	ng	99
32) Benzoic acid	9.48	122	151176	43.96	ng	84
33) 4-Chloroaniline	9.79	127	90902	10.82	ng	# 94
34) Hexachlorobutadiene	9.84	225	207125	48.83	ng	97
35) Caprolactam	10.42	113	96118m	54.02	ng	
36) 4-Chloro-3-methylphenol	10.64	107	400610	58.13	ng	92
37) 2-Methylnaphthalene	10.75	142	800075	56.15	ng	96
39) 1,2,4,5-Tetrachlorobenzene	11.03	216	347149	49.66	ng	99
40) Hexachlorocyclopentadiene	11.00	237	261631	89.66	ng	95

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	11.26	196	227806	56.46	nq	93
43) 2,4,5-Trichlorophenol	11.34	196	226292	55.07	nq	# 91
45) 1,1'-Biphenyl	11.52	154	972314	52.88	nq	97
46) 2-Chloronaphthalene	11.53	162	736861	52.39	nq	# 90
47) 2-Nitroaniline	11.75	65	303267	59.84	nq	# 74
48) Acenaphthylene	12.19	152	1192263	53.54	nq	99
49) Dimethylphthalate	12.08	163	953153	54.49	nq	100
50) 2,6-Dinitrotoluene	12.17	165	199740	56.67	nq	# 85
51) Acenaphthene	12.48	154	741723	54.19	nq	98
52) 3-Nitroaniline	12.43	138	107156	29.62	nq	# 83
53) 2,4-Dinitrophenol	12.63	184	156991	86.45	nq	# 74
54) Dibenzofuran	12.76	168	1078040	58.17	nq	95
55) 4-Nitrophenol	12.81	139	211368	122.48	nq	# 51
56) 2,4-Dinitrotoluene	12.83	165	279364	59.30	nq	90
57) Fluorene	13.31	166	894657	55.67	nq	99
58) 2,3,4,6-Tetrachlorophenol	13.00	232	211012	66.76	nq	97
59) Diethylphthalate	13.23	149	955135	56.16	nq	98
60) 4-Chlorophenyl-phenylether	13.35	204	411558	52.52	nq	91
61) 4-Nitroaniline	13.45	138	188187	55.74	nq	# 77
62) Azobenzene	13.60	77	1102380	62.48	nq	86
64) 4,6-Dinitro-2-methylphenol	13.50	198	129383	47.02	nq	# 69
65) n-Nitrosodiphenylamine	13.57	169	753110	53.81	nq	100
66) 4-Bromophenyl-phenylether	14.14	248	245599	51.88	nq	# 87
67) Hexachlorobenzene	14.21	284	253290	51.15	nq	# 83
68) Atrazine	14.48	200	241511	54.14	nq	93
69) Pentachlorophenol	14.56	266	210372	132.32	nq	96
70) Phenanthrene	14.87	178	1292295	53.91	nq	100
71) Anthracene	14.95	178	1282170	52.95	nq	100
72) Carbazole	15.23	167	1151021	55.73	nq	99
73) Di-n-butylphthalate	15.81	149	1463925	54.81	nq	# 98
74) Fluoranthene	16.62	202	1260717	52.45	nq	97
76) Benzidine	16.86	184	99105	11.21	nq	96
77) Pyrene	16.93	202	1317295	53.27	nq	99
79) Butylbenzylphthalate	17.83	149	605065	55.18	nq	# 67
80) Benzo(a)anthracene	18.50	228	1058289	54.16	nq	99
81) 3,3'-Dichlorobenzidine	18.49	252	166587	15.43	nq	# 96
82) Chrysene	18.55	228	1015821	52.38	nq	98
83) Bis(2-ethylhexyl)phthalate	18.57	149	823930	51.49	nq	# 95
84) Di-n-octyl phthalate	19.35	149	1273997	52.21	nq	# 87
85) Indeno(1,2,3-cd)pyrene	21.42	276	745116	47.93	nq	95
87) Benzo(b)fluoranthene	19.77	252	764963m	49.80	nq	
88) Benzo(k)fluoranthene	19.81	252	862399m	65.23	nq	
89) Benzo(a)pyrene	20.13	252	757193	56.99	nq	97
90) Dibenzo(a,h)anthracene	21.43	278	634405	55.99	nq	# 93
91) Benzo(g,h,i)perylene	21.77	276	625654	55.96	nq	94

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

