

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051616\
 Data File : BF087089.D
 Acq On : 16 May 2016 18:55
 Operator : UM/SJ
 Sample : H3074-01MS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HSR-WC-44-051216MS

Manual Integrations
 APPROVED

umangi
 5/17/2016 7:15:51 PM

Quant Time: May 17 13:25:46 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 17 01:12:04 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.60	152	128855	20.00	ng	0.00
21) Naphthalene-d8	7.90	136	501058	20.00	ng	0.00
38) Acenaphthene-d10	9.64	164	201505	20.00	ng	0.00
63) Phenanthrene-d10	11.12	188	321408	20.00	ng	0.00
75) Chrysene-d12	13.75	240	223730	20.00	ng	-0.01
86) Perylene-d12	15.11	264	209749	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.22	112	976386	128.33	ng	0.02
7) Phenol-d6	6.28	99	1122108	110.35	ng	0.00
23) Nitrobenzene-d5	7.19	82	919665	92.16	ng	0.00
41) 2,4,6-Tribromophenol	10.43	330	250075	117.23	ng	-0.01
44) 2-Fluorobiphenyl	8.98	172	1368548	114.14	ng	0.00
78) Terphenyl-d14	12.71	244	1003370	95.15	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.23	88	125174	33.51	ng	# 67
3) Pyridine	2.88	79	209321m	22.92	ng	
4) n-Nitrosodimethylamine	2.78	42	288817	43.61	ng	# 47
6) Aniline	6.28	93	112602	9.16	ng	# 1
8) 2-Chlorophenol	6.40	128	364833	39.74	ng	# 78
9) Benzaldehyde	6.16	77	27991	4.75	ng	97
10) Phenol	6.30	94	438454	36.28	ng	75
11) bis(2-Chloroethyl)ether	6.35	93	376352	39.93	ng	84
12) 1,3-Dichlorobenzene	6.55	146	388546	39.11	ng	97
13) 1,4-Dichlorobenzene	6.63	146	399200	39.40	ng	98
14) 1,2-Dichlorobenzene	6.78	146	341787	38.70	ng	95
15) Benzyl Alcohol	6.78	79	342455	48.77	ng	92
16) 2,2'-oxybis(1-Chloropropan	6.90	45	800826	39.08	ng	70
17) 2-Methylphenol	6.90	107	302789m	41.45	ng	
18) Hexachloroethane	7.12	117	118180	31.00	ng	# 81
19) n-Nitroso-di-n-propylamine	7.04	70	288076m	40.54	ng	
20) 3+4-Methylphenols	7.06	107	410989	43.07	ng	# 60
22) Acetophenone	7.03	105	536769	45.35	ng	# 94
24) Nitrobenzene	7.20	77	407409	39.69	ng	92
25) Isophorone	7.45	82	784018	42.35	ng	94
26) 2-Nitrophenol	7.52	139	161451	35.78	ng	# 72
27) 2,4-Dimethylphenol	7.58	122	317490	41.31	ng	# 85
28) bis(2-Chloroethoxy)methane	7.67	93	461969	46.26	ng	98
29) 2,4-Dichlorophenol	7.78	162	287783	42.54	ng	94
30) 1,2,4-Trichlorobenzene	7.84	180	300643	39.75	ng	95
31) Naphthalene	7.92	128	1075309	42.85	ng	99
32) Benzoic acid	7.74	122	171425	37.97	ng	# 80
33) 4-Chloroaniline	7.99	127	75903	7.78	ng	# 92
34) Hexachlorobutadiene	8.03	225	170866	38.93	ng	99
35) Caprolactam	8.36	113	72028	31.29	ng	# 46
36) 4-Chloro-3-methylphenol	8.49	107	322213	39.27	ng	88
37) 2-Methylnaphthalene	8.60	142	645104	43.97	ng	# 97
39) 1,2,4,5-Tetrachlorobenzene	8.78	216	274495	46.85	ng	99
40) Hexachlorocyclopentadiene	8.75	237	37398	13.45	ng	93

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.90	196	182541	46.59	nq	96
43) 2,4,5-Trichlorophenol	8.96	196	167614	41.37	nq	95
45) 1,1'-Biphenyl	9.07	154	740668	46.21	nq	96
46) 2-Chloronaphthalene	9.10	162	549592	43.77	nq	93
47) 2-Nitroaniline	9.21	65	261307	56.94	nq	# 83
48) Acenaphthylene	9.51	152	844037	45.70	nq	99
49) Dimethylphthalate	9.38	163	685443	44.59	nq	99
50) 2,6-Dinitrotoluene	9.45	165	140308	43.37	nq	95
51) Acenaphthene	9.68	154	548413	47.73	nq	98
52) 3-Nitroaniline	9.62	138	78273	22.98	nq	97
53) 2,4-Dinitrophenol	9.74	184	21788	19.60	nq	# 74
54) Dibenzofuran	9.85	168	740885	50.23	nq	95
55) 4-Nitrophenol	9.82	139	119265m	54.08	nq	
56) 2,4-Dinitrotoluene	9.85	165	138058	34.48	nq	# 37
57) Fluorene	10.19	166	611390	48.61	nq	99
58) 2,3,4,6-Tetrachlorophenol	9.99	232	133771	43.86	nq	97
59) Diethylphthalate	10.08	149	643193	42.93	nq	99
60) 4-Chlorophenyl-phenylether	10.19	204	264572	49.58	nq	# 76
61) 4-Nitroaniline	10.24	138	149267	45.63	nq	89
62) Azobenzene	10.34	77	703414	43.52	nq	82
64) 4,6-Dinitro-2-methylphenol	10.26	198	18624	9.33	nq	# 23
65) n-Nitrosodiphenylamine	10.31	169	486685	49.29	nq	100
66) 4-Bromophenyl-phenylether	10.67	248	159668	49.00	nq	97
67) Hexachlorobenzene	10.74	284	172759	45.36	nq	99
68) Atrazine	10.84	200	131145	39.76	nq	95
69) Pentachlorophenol	10.95	266	148054	84.12	nq	98
70) Phenanthrene	11.14	178	753921	43.95	nq	100
71) Anthracene	11.20	178	817774	46.62	nq	99
72) Carbazole	11.36	167	669040	44.37	nq	98
73) Di-n-butylphthalate	11.70	149	952514	51.72	nq	98
74) Fluoranthene	12.33	202	808325	45.82	nq	94
76) Benzidine	12.47	184	225692	39.57	nq	96
77) Pyrene	12.56	202	789961	46.12	nq	99
79) Butylbenzylphthalate	13.19	149	364430	50.30	nq	90
80) Benzo(a)anthracene	13.74	228	622918	49.60	nq	100
81) 3,3'-Dichlorobenzidine	13.71	252	201171	25.21	nq	# 97
82) Chrysene	13.78	228	632448	49.50	nq	99
83) Bis(2-ethylhexyl)phthalate	13.75	149	490722	56.31	nq	98
84) Di-n-octyl phthalate	14.35	149	829382	57.45	nq	# 83
85) Indeno(1,2,3-cd)pyrene	16.37	276	522441	49.15	nq	96
87) Benzo(b)fluoranthene	14.74	252	568623m	41.71	nq	
88) Benzo(k)fluoranthene	14.77	252	538706m	48.26	nq	
89) Benzo(a)pyrene	15.06	252	544432	46.30	nq	97
90) Dibenzo(a,h)anthracene	16.38	278	450196	45.43	nq	# 93
91) Benzo(g,h,i)perylene	16.73	276	420021	41.71	nq	95

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

