

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061716\
 Data File : BF088093.D
 Acq On : 17 Jun 2016 15:15
 Operator : UM/SJ
 Sample : H3510-01MS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HSR-WC-52-060916MS

Manual Integrations
 APPROVED

sohil
 6/20/2016 6:50:49 PM

Quant Time: Jun 17 17:57:29 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 17 16:09:13 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	117684	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	503534	20.00	ng	0.00
38) Acenaphthene-d10	9.74	164	231514	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	428806	20.00	ng	0.00
75) Chrysene-d12	13.85	240	280792	20.00	ng	0.00
86) Perylene-d12	15.23	264	45762	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.31	112	758680	110.21	ng	0.03
7) Phenol-d6	6.35	99	857852	102.83	ng	0.00
23) Nitrobenzene-d5	7.27	82	608422	70.56	ng	-0.01
41) 2,4,6-Tribromophenol	10.52	330	237368	97.10	ng	0.00
44) 2-Fluorobiphenyl	9.06	172	1205177	81.55	ng	-0.01
78) Terphenyl-d14	12.80	244	942983	76.42	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.40	88	100549	30.06	ng	# 87
3) Pyridine	3.10	79	129670m	16.42	ng	
4) n-Nitrosodimethylamine	2.96	42	118875	35.54	ng	82
8) 2-Chlorophenol	6.49	128	323912	39.75	ng	88
9) Benzaldehyde	6.25	77	15575	2.68	ng	98
10) Phenol	6.36	94	267224	30.20	ng	86
11) bis(2-Chloroethyl)ether	6.44	93	288089	39.38	ng	89
12) 1,3-Dichlorobenzene	6.64	146	333911m	38.19	ng	
13) 1,4-Dichlorobenzene	6.72	146	334791	38.02	ng	96
14) 1,2-Dichlorobenzene	6.87	146	313354m	39.13	ng	
15) Benzyl Alcohol	6.86	79	68327	10.47	ng	91
16) 2,2'-oxybis(1-Chloropropan	6.98	45	375576	38.00	ng	60
17) 2-Methylphenol	6.97	107	242037	36.63	ng	98
18) Hexachloroethane	7.21	117	125277	40.09	ng	100
19) n-Nitroso-di-n-propylamine	7.13	70	216260	40.52	ng	# 83
20) 3+4-Methylphenols	7.13	107	301943	39.16	ng	# 80
22) Acetophenone	7.12	105	444454	39.50	ng	98
24) Nitrobenzene	7.29	77	351588	42.09	ng	99
25) Isophorone	7.53	82	611120	38.26	ng	100
26) 2-Nitrophenol	7.60	139	182378	40.76	ng	89
27) 2,4-Dimethylphenol	7.66	122	318490	38.66	ng	97
28) bis(2-Chloroethoxy)methane	7.75	93	279524	28.22	ng	99
29) 2,4-Dichlorophenol	7.85	162	282700	41.10	ng	98
30) 1,2,4-Trichlorobenzene	7.93	180	285498	38.12	ng	98
31) Naphthalene	8.00	128	995210	39.94	ng	99
32) Benzoic acid	7.80	122	208194	45.89	ng	93
34) Hexachlorobutadiene	8.12	225	145884	36.93	ng	99
35) Caprolactam	8.46	113	81136m	35.61	ng	
36) 4-Chloro-3-methylphenol	8.56	107	292618	39.78	ng	97
37) 2-Methylnaphthalene	8.70	142	626421	38.14	ng	# 97
39) 1,2,4,5-Tetrachlorobenzene	8.87	216	274734	45.91	ng	# 1
40) Hexachlorocyclopentadiene	8.86	237	193471	51.81	ng	98
42) 2,4,6-Trichlorophenol	8.98	196	191531	41.37	ng	96
43) 2,4,5-Trichlorophenol	9.04	196	183801	38.16	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
45) 1,1'-Biphenyl	9.16	154	774970	44.53	nq	98
46) 2-Chloronaphthalene	9.19	162	605289	40.40	nq	98
47) 2-Nitroaniline	9.29	65	104759	23.70	nq	# 69
48) Acenaphthylene	9.60	152	773868	32.47	nq	100
49) Dimethylphthalate	9.47	163	703217	41.93	nq	99
50) 2,6-Dinitrotoluene	9.53	165	172383	44.88	nq	# 64
51) Acenaphthene	9.77	154	531609	35.76	nq	99
53) 2,4-Dinitrophenol	9.82	184	160404	76.96	nq	# 84
54) Dibenzofuran	9.94	168	736545	35.98	nq	# 88
55) 4-Nitrophenol	9.88	139	201933	58.71	nq	# 79
56) 2,4-Dinitrotoluene	9.94	165	207173	41.97	nq	# 69
57) Fluorene	10.28	166	571802	40.89	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.07	232	148844	39.32	nq	# 53
59) Diethylphthalate	10.17	149	692240	40.78	nq	99
60) 4-Chlorophenyl-phenylether	10.28	204	261619	38.43	nq	# 84
61) 4-Nitroaniline	10.31	138	46969	11.17	nq	97
62) Azobenzene	10.43	77	590034	35.15	nq	95
64) 4,6-Dinitro-2-methylphenol	10.34	198	104691	39.51	nq	95
65) n-Nitrosodiphenylamine	10.40	169	77309	5.43	nq	99
66) 4-Bromophenyl-phenylether	10.76	248	173121	37.63	nq	# 84
67) Hexachlorobenzene	10.83	284	198053	41.43	nq	# 80
69) Pentachlorophenol	11.03	266	182649	72.49	nq	97
70) Phenanthrene	11.24	178	964739	42.87	nq	99
71) Anthracene	11.29	178	820834	36.16	nq	100
72) Carbazole	11.45	167	129163	6.08	nq	96
73) Di-n-butylphthalate	11.78	149	1043022	38.79	nq	99
74) Fluoranthene	12.42	202	803735	33.85	nq	98
77) Pvrene	12.65	202	759557	36.45	nq	100
79) Butylbenzylphthalate	13.28	149	437246	47.46	nq	99
80) Benzo(a)anthracene	13.84	228	620916	38.80	nq	100
82) Chrysene	13.87	228	568008	37.73	nq	99
83) Bis(2-ethylhexyl)phthalate	13.84	149	492504	43.92	nq	# 100
84) Di-n-octyl phthalate	14.45	149	931079	51.10	nq	# 100
85) Indeno(1,2,3-cd)pyrene	16.54	276	371889	26.59	nq	# 100
87) Benzo(b)fluoranthene	14.85	252	567990m	178.08	nq	
88) Benzo(k)fluoranthene	14.88	252	538252m	223.71	nq	
89) Benzo(a)pyrene	15.18	252	347421	134.63	nq	# 96
90) Dibenzo(a,h)anthracene	16.55	278	439939	183.56	nq	99
91) Benzo(g,h,i)perylene	16.93	276	291345	118.17	nq	99

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

