

Data Path : Z:\HPCHEM1\BNA F\DATA\BF010715\
 Data File : BF076771.D
 Acq On : 7 Jan 2015 19:15
 Operator : IZ / TP
 Sample : G1022-03MS
 Misc : W
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-H2O-FP3-65MS

Manual Integrations
 APPROVED

tejaskumar
 1/8/2015 3:46:17 PM

Quant Time: Jan 08 09:27:23 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121714.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 08 09:08:31 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.14	152	43491	20.00	ng	0.00
21) Naphthalene-d8	11.04	136	188842	20.00	ng	0.00
38) Acenaphthene-d10	15.19	164	104835	20.00	ng	0.00
63) Phenanthrene-d10	17.90	188	168777	20.00	ng	0.00
75) Chrysene-d12	21.39	240	143260	20.00	ng	-0.01
86) Perylene-d12	23.07	264	61254	20.00	ng	-0.05

System Monitoring Compounds

5) 2-Fluorophenol	5.21	112	300784	117.25	ng	0.00
7) Phenol-d6	7.59	99	364425	116.31	ng	0.07
23) Nitrobenzene-d5	9.44	82	255379	80.43	ng	0.01
41) 2,4,6-Tribromophenol	16.83	330	108248	122.17	ng	0.01
44) 2-Fluorobiphenyl	13.71	172	528620	78.47	ng	0.00
78) Terphenyl-d14	20.12	244	540144	83.17	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.40	88	33575	30.62	ng	# 100
3) Pyridine	1.93	79	82403m	29.85	ng	
4) n-Nitrosodimethylamine	1.86	42	43017	32.21	ng	90
8) 2-Chlorophenol	7.67	128	115933	39.46	ng	98
9) Benzaldehyde	7.14	77	109747	46.04	ng	96
10) Phenol	7.62	94	138473m	36.42	ng	
11) bis(2-Chloroethyl)ether	7.66	93	112464	41.09	ng	91
12) 1,3-Dichlorobenzene	7.98	146	117590	34.48	ng	96
13) 1,4-Dichlorobenzene	8.17	146	119600	34.62	ng	95
14) 1,2-Dichlorobenzene	8.49	146	116795	35.65	ng	95
15) Benzyl Alcohol	8.66	79	1671839	616.22	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.92	45	137841	34.76	ng	# 36
17) 2-Methylphenol	8.93	107	91074	37.73	ng	# 91
18) Hexachloroethane	9.25	117	45329	36.72	ng	# 88
19) n-Nitroso-di-n-propylamine	9.20	70	93417m	40.83	ng	
20) 3+4-Methylphenols	9.29	107	119894	36.78	ng	95
22) Acetophenone	9.13	105	177985	39.32	ng	# 96
24) Nitrobenzene	9.48	77	125389	38.08	ng	94
25) Isophorone	10.08	82	240943	39.37	ng	94
26) 2-Nitrophenol	10.22	139	60974	37.92	ng	89
27) 2,4-Dimethylphenol	10.49	122	102975	39.57	ng	96
28) bis(2-Chloroethoxy)methane	10.69	93	110070	30.64	ng	97
29) 2,4-Dichlorophenol	10.82	162	100708	39.37	ng	95
30) 1,2,4-Trichlorobenzene	10.95	180	100267	36.36	ng	# 94
31) Naphthalene	11.10	128	340280	36.37	ng	99
32) Benzoic acid	11.80	122	3131680m	1798.94	ng	
34) Hexachlorobutadiene	11.45	225	57558	35.90	ng	91
36) 4-Chloro-3-methylphenol	12.63	107	106747	37.50	ng	92
37) 2-Methylnaphthalene	12.76	142	241213	36.96	ng	96
39) 1,2,4,5-Tetrachlorobenzene	13.16	216	93870	36.40	ng	# 76
40) Hexachlorocyclopentadiene	13.15	237	113311	76.25	ng	94
42) 2,4,6-Trichlorophenol	13.50	196	69755	37.28	ng	98
43) 2,4,5-Trichlorophenol	13.59	196	72891	36.21	ng	# 92
45) 1,1'-Biphenyl	13.91	154	291046	37.37	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
46) 2-Chloronaphthalene	13.89	162	227948	36.90	nq	97
47) 2-Nitroaniline	14.22	65	51797	27.57	nq	97
48) Acenaphthylene	14.84	152	366798	34.91	nq	98
49) Dimethylphthalate	14.79	163	678333	91.47	nq	100
50) 2,6-Dinitrotoluene	14.87	165	59462	39.13	nq	# 67
51) Acenaphthene	15.27	154	212097	35.68	nq	98
53) 2,4-Dinitrophenol	15.50	184	52315	65.10	nq	# 89
54) Dibenzofuran	15.68	168	331771	38.59	nq	99
55) 4-Nitrophenol	15.80	139	81477	61.60	nq	# 81
56) 2,4-Dinitrotoluene	15.79	165	84022	41.49	nq	# 62
57) Fluorene	16.38	166	269300	37.29	nq	97
58) 2,3,4,6-Tetrachlorophenol	16.00	232	55875	37.32	nq	# 100
59) Diethylphthalate	16.36	149	1317146	173.89	nq	98
60) 4-Chlorophenyl-phenylether	16.46	204	119785	38.33	nq	92
61) 4-Nitroaniline	16.51	138	12686	7.00	nq	85
62) Azobenzene	16.72	77	255106	34.18	nq	87
64) 4,6-Dinitro-2-methylphenol	16.56	198	38141	36.13	nq	# 79
65) n-Nitrosodiphenylamine	16.69	169	90165	15.21	nq	99
66) 4-Bromophenyl-phenylether	17.28	248	68059	41.40	nq	# 78
67) Hexachlorobenzene	17.31	284	75005	39.09	nq	92
68) Atrazine	17.66	200	43749	24.70	nq	79
69) Pentachlorophenol	17.65	266	77471	82.48	nq	98
70) Phenanthrene	17.93	178	384172	37.73	nq	99
71) Anthracene	18.01	178	382821	38.32	nq	97
72) Carbazole	18.29	167	104388	10.77	nq	97
73) Di-n-butylphthalate	18.87	149	575791	48.38	nq	99
74) Fluoranthene	19.57	202	380187	36.56	nq	97
77) Pvrene	19.85	202	355768	35.50	nq	98
79) Butylbenzylphthalate	20.78	149	321796	67.16	nq	99
80) Benzo(a)anthracene	21.37	228	323554	36.27	nq	98
82) Chrysene	21.42	228	315282	38.62	nq	99
83) Bis(2-ethylhexyl)phthalate	21.51	149	366221	50.52	nq	# 96
84) Di-n-octyl phthalate	22.28	149	523625	42.31	nq	# 99
87) Benzo(b)fluoranthene	22.64	252	310467m	76.83	nq	
88) Benzo(k)fluoranthene	22.68	252	275751m	73.27	nq	
89) Benzo(a)pvrene	23.00	252	201400	55.85	nq	# 97
90) Dibenzo(a,h)anthracene	24.24	278	247944	69.38	nq	99
91) Benzo(g,h,i)perylene	24.53	276	206178	59.29	nq	96

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

