

Data Path : Z:\HPCHEM1\BNA F\DATA\BF123114\
 Data File : BF076705.D
 Acq On : 31 Dec 2014 16:59
 Operator : IZ / TP
 Sample : IDOC-02-W-RS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 IDOC-02-W-RS

Quant Time: Jan 01 06:27:49 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121714.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 31 23:59:08 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	40814	20.00	ng	0.00
21) Naphthalene-d8	8.30	136	170063	20.00	ng	-0.01
38) Acenaphthene-d10	10.45	164	98903	20.00	ng	0.00
63) Phenanthrene-d10	12.25	188	164054	20.00	ng	-0.01
75) Chrysene-d12	15.50	240	149474	20.00	ng	0.00
86) Perylene-d12	17.16	264	136975	20.00	ng	-0.03

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	4.90	112	146703	60.94	ng	0.00
7) Phenol-d6	6.32	99	109603	37.28	ng	0.00
23) Nitrobenzene-d5	7.43	82	278599	97.43	ng	0.00
41) 2,4,6-Tribromophenol	11.42	330	118709	142.01	ng	0.00
44) 2-Fluorobiphenyl	9.66	172	559170	87.98	ng	0.00
78) Terphenyl-d14	14.24	244	605352	89.33	ng	-0.01

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	1.41	88	13837	13.45	ng	# 100
3) Pyridine	1.94	79	32123	12.40	ng	# 82
4) n-Nitrosodimethylamine	1.88	42	20597	16.44	ng	# 83
6) Aniline	6.30	93	64082m	15.26	ng	
8) 2-Chlorophenol	6.45	128	100018	36.28	ng	95
9) Benzaldehyde	6.15	77	9287	4.15	ng	97
10) Phenol	6.33	94	43292	12.13	ng	83
11) bis(2-Chloroethyl)ether	6.42	93	126806	49.37	ng	94
12) 1,3-Dichlorobenzene	6.63	146	128105	40.02	ng	# 92
13) 1,4-Dichlorobenzene	6.74	146	133827	41.28	ng	96
14) 1,2-Dichlorobenzene	6.93	146	128471	41.79	ng	97
15) Benzyl Alcohol	6.93	79	73243	28.77	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.12	45	180541	48.52	ng	98
17) 2-Methylphenol	7.10	107	67597	29.84	ng	95
18) Hexachloroethane	7.34	117	48952	42.25	ng	97
19) n-Nitroso-di-n-propylamine	7.27	70	93390	43.49	ng	96
20) 3+4-Methylphenols	7.30	107	76758	25.09	ng	93
22) Acetophenone	7.25	105	215085	52.77	ng	# 98
24) Nitrobenzene	7.45	77	151401	51.06	ng	99
25) Isophorone	7.76	82	278515	50.53	ng	94
26) 2-Nitrophenol	7.85	139	66732	46.09	ng	99
27) 2,4-Dimethylphenol	7.94	122	99124	42.29	ng	93
28) bis(2-Chloroethoxy)methane	8.06	93	169188	52.29	ng	98
29) 2,4-Dichlorophenol	8.15	162	101184	43.92	ng	93
30) 1,2,4-Trichlorobenzene	8.24	180	112121	45.15	ng	# 93
31) Naphthalene	8.33	128	398229	47.27	ng	98
32) Benzoic acid	8.06	122	16705	15.56	ng	99
33) 4-Chloroaniline	8.42	127	83796	24.37	ng	96
34) Hexachlorobutadiene	8.49	225	65053	45.06	ng	93
35) Caprolactam	8.85	113	4618	6.95	ng	# 76
36) 4-Chloro-3-methylphenol	9.04	107	98399	38.39	ng	# 86
37) 2-Methylnaphthalene	9.19	142	269281	45.81	ng	97
39) 1,2,4,5-Tetrachlorobenzene	9.38	216	114712	47.15	ng	# 79
40) Hexachlorocyclopentadiene	9.37	237	104175	74.48	ng	94

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.54	196	76285	43.22	nq	98
43) 2,4,5-Trichlorophenol	9.59	196	80623	42.45	nq	# 91
45) 1,1'-Biphenyl	9.77	154	339546	46.21	nq	98
46) 2-Chloronaphthalene	9.77	162	264912	45.46	nq	96
47) 2-Nitroaniline	9.92	65	93884	52.96	nq	# 86
48) Acenaphthylene	10.28	152	436668	44.06	nq	99
49) Dimethylphthalate	10.17	163	342753	48.99	nq	100
50) 2,6-Dinitrotoluene	10.24	165	69605	48.56	nq	95
51) Acenaphthene	10.49	154	252330	45.00	nq	98
52) 3-Nitroaniline	10.42	138	53279	32.73	nq	94
53) 2,4-Dinitrophenol	10.56	184	54734	70.06	nq	# 60
54) Dibenzofuran	10.70	168	370599	45.69	nq	96
55) 4-Nitrophenol	10.68	139	34404m	27.57	nq	
56) 2,4-Dinitrotoluene	10.72	165	98525	51.57	nq	# 78
57) Fluorene	11.12	166	321686	47.21	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.87	232	63945	45.28	nq	# 100
59) Diethylphthalate	11.04	149	362957	50.79	nq	98
60) 4-Chlorophenyl-phenylether	11.14	204	144843	49.12	nq	# 86
61) 4-Nitroaniline	11.17	138	61534	35.97	nq	86
62) Azobenzene	11.34	77	343753	48.82	nq	95
64) 4,6-Dinitro-2-methylphenol	11.21	198	41475	40.41	nq	98
65) n-Nitrosodiphenylamine	11.29	169	251540	43.65	nq	97
66) 4-Bromophenyl-phenylether	11.74	248	82909	51.88	nq	# 83
67) Hexachlorobenzene	11.78	284	90444	48.49	nq	97
68) Atrazine	11.98	200	87737	50.95	nq	99
69) Pentachlorophenol	12.04	266	88442	96.87	nq	96
70) Phenanthrene	12.29	178	460659	46.54	nq	99
71) Anthracene	12.36	178	474292	48.84	nq	98
72) Carbazole	12.56	167	416873	44.26	nq	97
73) Di-n-butylphthalate	13.05	149	598664	51.75	nq	# 97
74) Fluoranthene	13.74	202	486554	48.14	nq	97
76) Benzidine	13.95	184	177182	28.44	nq	99
77) Pyrene	14.01	202	521054	49.83	nq	99
79) Butylbenzylphthalate	14.88	149	267060	53.42	nq	99
80) Benzo(a)anthracene	15.49	228	450867	48.44	nq	98
81) 3,3'-Dichlorobenzidine	15.49	252	113076	36.34	nq	99
82) Chrysene	15.52	228	428337	50.29	nq	99
83) Bis(2-ethylhexyl)phthalate	15.60	149	418558	55.34	nq	# 98
84) Di-n-octyl phthalate	16.38	149	707983	54.82	nq	100
85) Indeno(1,2,3-cd)pyrene	18.32	276	450127	48.16	nq	96
87) Benzo(b)fluoranthene	16.75	252	443822	49.12	nq	97
88) Benzo(k)fluoranthene	16.77	252	389215m	46.25	nq	
89) Benzo(a)pyrene	17.10	252	392816	48.72	nq	97
90) Dibenzo(a,h)anthracene	18.35	278	387677	48.51	nq	99
91) Benzo(g,h,i)perylene	18.63	276	389412	50.08	nq	98

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

