

Data Path : Z:\HPCHEM1\BNA F\DATA\BF111414\
 Data File : BF075512.D
 Acq On : 14 Nov 2014 23:52
 Operator : TP / JJ
 Sample : PB80327BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB80327BS

Manual Integrations
 APPROVED

mohammad
 11/17/2014 7:34:41 PM

Quant Time: Nov 16 04:25:14 2014
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF111214.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 12 18:46:55 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	93331	20.00	ng	-0.02
21) Naphthalene-d8	9.13	136	397646	20.00	ng	-0.02
38) Acenaphthene-d10	11.32	164	203868	20.00	ng	-0.02
63) Phenanthrene-d10	13.16	188	326652	20.00	ng	-0.03
75) Chrysene-d12	16.46	240	359376	20.00	ng	-0.08
86) Perylene-d12	18.20	264	384215	20.00	ng	-0.24

System Monitoring Compounds

5) 2-Fluorophenol	5.84	112	426379	80.71	ng	-0.02
7) Phenol-d6	7.10	99	525423	82.70	ng	-0.01
23) Nitrobenzene-d5	8.24	82	283825	52.33	ng	-0.02
41) 2,4,6-Tribromophenol	12.30	330	114371	66.98	ng	-0.02
44) 2-Fluorobiphenyl	10.48	172	544904	48.30	ng	-0.01
78) Terphenyl-d14	15.16	244	560033	41.95	ng	-0.03

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.48	88	49792	22.45	ng	97
3) Pyridine	3.27	79	141120m	23.68	ng	
4) n-Nitrosodimethylamine	3.19	42	76371m	24.21	ng	
6) Aniline	7.13	93	118320	12.89	ng	98
8) 2-Chlorophenol	7.28	128	143582	23.48	ng	95
10) Phenol	7.11	94	198077	25.56	ng	87
11) bis(2-Chloroethyl)ether	7.24	93	150473	25.63	ng	# 81
12) 1,3-Dichlorobenzene	7.48	146	153077	22.97	ng	97
13) 1,4-Dichlorobenzene	7.57	146	150741	22.24	ng	99
14) 1,2-Dichlorobenzene	7.76	146	145018	22.82	ng	# 91
15) Benzyl Alcohol	7.73	79	118143	23.69	ng	95
16) 2,2'-oxybis(1-Chloropropan	7.91	45	285753	23.51	ng	95
17) 2-Methylphenol	7.86	107	117470	23.20	ng	98
18) Hexachloroethane	8.17	117	54295	23.36	ng	98
19) n-Nitroso-di-n-propylamine	8.06	70	100287	23.17	ng	94
20) 3+4-Methylphenols	8.06	107	154599	23.32	ng	90
22) Acetophenone	8.06	105	207499	24.19	ng	# 96
24) Nitrobenzene	8.26	77	155801	24.79	ng	99
25) Isophorone	8.56	82	285728	23.07	ng	98
26) 2-Nitrophenol	8.65	139	65161	22.75	ng	95
27) 2,4-Dimethylphenol	8.71	122	129971	23.41	ng	99
28) bis(2-Chloroethoxy)methane	8.84	93	186006	24.65	ng	98
29) 2,4-Dichlorophenol	8.95	162	112665	23.17	ng	97
30) 1,2,4-Trichlorobenzene	9.06	180	111107	21.76	ng	97
31) Naphthalene	9.16	128	427473	24.11	ng	99
32) Benzoic acid	8.80	122	90125m	26.87	ng	
33) 4-Chloroaniline	9.22	127	102651	13.33	ng	98
34) Hexachlorobutadiene	9.32	225	57668	20.90	ng	96
35) Caprolactam	9.64	113	36556	22.68	ng	92
36) 4-Chloro-3-methylphenol	9.82	107	114184	21.40	ng	91
37) 2-Methylnaphthalene	10.01	142	283084	23.40	ng	98
39) 1,2,4,5-Tetrachlorobenzene	10.22	216	99357	23.71	ng	98
40) Hexachlorocyclopentadiene	10.21	237	128190	50.36	ng	100
42) 2,4,6-Trichlorophenol	10.37	196	73810	22.15	ng	99

Data Path : Z:\HPCHEM1\BNA F\DATA\BF111414\
 Data File : BF075512.D
 Acq On : 14 Nov 2014 23:52
 Operator : TP / JJ
 Sample : PB80327BS
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleID :
 PB80327BS

Manual Integrations
 APPROVED

mohammad
 11/17/2014 7:34:41 PM

Quant Time: Nov 16 04:25:14 2014
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF111214.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 12 18:46:55 2014
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
43) 2,4,5-Trichlorophenol	10.41	196	73746	21.25	nq	96
45) 1,1'-Biphenyl	10.60	154	322217	23.43	nq	99
46) 2-Chloronaphthalene	10.62	162	257916	23.99	nq	98
47) 2-Nitroaniline	10.74	65	78760	24.35	nq	87
48) Acenaphthylene	11.13	152	443906	25.04	nq	99
49) Dimethylphthalate	10.98	163	292296	20.97	nq	99
50) 2,6-Dinitrotoluene	11.05	165	63509	23.14	nq	# 81
51) Acenaphthene	11.35	154	267110	24.96	nq	99
52) 3-Nitroaniline	11.26	138	58008	17.92	nq	# 87
53) 2,4-Dinitrophenol	11.38	184	50880	41.90	nq	86
54) Dibenzofuran	11.57	168	353599	23.12	nq	94
55) 4-Nitrophenol	11.45	139	119208	46.07	nq	95
56) 2,4-Dinitrotoluene	11.56	165	83860	23.25	nq	94
57) Fluorene	11.99	166	286168	23.14	nq	99
58) 2,3,4,6-Tetrachlorophenol	11.72	232	64099	23.27	nq	# 90
59) Diethylphthalate	11.86	149	286717	19.83	nq	96
60) 4-Chlorophenyl-phenylether	12.00	204	119070	22.39	nq	# 86
61) 4-Nitroaniline	12.01	138	73696	22.78	nq	84
62) Azobenzene	12.20	77	289066	22.72	nq	88
64) 4,6-Dinitro-2-methylphenol	12.05	198	33301	20.16	nq	# 41
65) n-Nitrosodiphenylamine	12.14	169	234668	22.71	nq	100
66) 4-Bromophenyl-phenylether	12.61	248	72041	23.44	nq	# 85
67) Hexachlorobenzene	12.68	284	82240	23.56	nq	# 83
68) Atrazine	12.81	200	69517	21.68	nq	96
69) Pentachlorophenol	12.92	266	96847	44.21	nq	98
70) Phenanthrene	13.19	178	430891	25.06	nq	99
71) Anthracene	13.26	178	455800	25.65	nq	99
72) Carbazole	13.45	167	412480	23.72	nq	99
73) Di-n-butylphthalate	13.90	149	480013	19.99	nq	99
74) Fluoranthene	14.68	202	440320	23.91	nq	92
76) Benzidine	14.84	184	233805m	20.46	nq	
77) Pyrene	14.95	202	465333m	22.27	nq	
79) Butylbenzylphthalate	15.76	149	211408	18.02	nq	# 77
80) Benzo(a)anthracene	16.44	228	427395	22.64	nq	99
81) 3,3'-Dichlorobenzidine	16.41	252	69075	9.91	nq	# 96
82) Chrysene	16.48	228	402061	24.62	nq	99
83) Bis(2-ethylhexyl)phthalate	16.48	149	310202	16.78	nq	100
84) Di-n-octyl phthalate	17.26	149	573938m	17.64	nq	
85) Indeno(1,2,3-cd)pyrene	19.75	276	554428m	27.99	nq	
87) Benzo(b)fluoranthene	17.73	252	522115	25.42	nq	99
88) Benzo(k)fluoranthene	17.76	252	381433m	20.69	nq	
89) Benzo(a)pyrene	18.13	252	447757	24.36	nq	# 95
90) Dibenzo(a,h)anthracene	19.77	278	468555	24.41	nq	# 94
91) Benzo(g,h,i)perylene	20.22	276	473922	25.91	nq	96

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

Instrument :
BNA_F
ClientSampleId :
PB80327BS

mohammad
11/17/2014 7:34:41 PM

[illegible]