

Data Path : Z:\HPCHEM1\BNA G\DATA\BG012115\
 Data File : BG015909.D
 Acq On : 21 Jan 2015 19:17
 Operator : TP/IZ
 Sample : G1100-18DL 10X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 8024DL

Quant Time: Jan 22 00:57:32 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 22 00:16:08 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	94937	20.00	ng	0.00
21) Naphthalene-d8	10.47	136	393124	20.00	ng	0.00
38) Acenaphthene-d10	14.33	164	341201	20.00	ng	0.00
63) Phenanthrene-d10	17.08	188	840794	20.00	ng	0.00
75) Chrysene-d12	21.27	240	1142091	20.00	ng	0.00
86) Perylene-d12	23.52	264	1083735	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.28	112	83943	13.17	ng	0.00
7) Phenol-d6	6.87	99	145476	14.16	ng	0.00
23) Nitrobenzene-d5	8.85	82	125481	9.20	ng	0.00
41) 2,4,6-Tribromophenol	15.83	330	75894	12.15	ng	0.00
44) 2-Fluorobiphenyl	12.95	172	213563	9.82	ng	0.00
78) Terphenyl-d14	19.71	244	540814	8.73	ng	0.00

Target Compounds

						Qvalue
10) Phenol	6.89	94	214701	18.34	ng	97
11) bis(2-Chloroethyl)ether	7.11	93	183142	20.24	ng	94
12) 1,3-Dichlorobenzene	7.58	146	155968m	21.54	ng	
13) 1,4-Dichlorobenzene	7.72	146	174278	23.32	ng	97
14) 1,2-Dichlorobenzene	8.03	146	107741	14.71	ng	# 85
26) 2-Nitrophenol	9.59	139	23628	6.65	ng	96
28) bis(2-Chloroethoxy)methane	9.89	93	171298	14.67	ng	# 93
30) 1,2,4-Trichlorobenzene	10.34	180	221724	25.76	ng	99
31) Naphthalene	10.52	128	302099	15.01	ng	97
34) Hexachlorobutadiene	10.80	225	125497	14.91	ng	93
36) 4-Chloro-3-methylphenol	11.77	107	241430	23.12	ng	93
37) 2-Methylnaphthalene	12.14	142	352194	22.04	ng	89
42) 2,4,6-Trichlorophenol	12.76	196	213850	28.06	ng	94
43) 2,4,5-Trichlorophenol	12.83	196	77526	9.00	ng	95
46) 2-Chloronaphthalene	13.21	162	87299	4.80	ng	# 77
48) Acenaphthylene	14.05	152	770507	27.90	ng	# 96
49) Dimethylphthalate	13.79	163	249886	10.18	ng	# 93
50) 2,6-Dinitrotoluene	13.92	165	160026	28.87	ng	98
51) Acenaphthene	14.39	154	355511	20.49	ng	93
55) 4-Nitrophenol	14.57	139	30380	8.47	ng	# 88
56) 2,4-Dinitrotoluene	14.70	165	142838	17.61	ng	# 92
57) Fluorene	15.38	166	268532	10.40	ng	93
59) Diethylphthalate	15.16	149	404025	16.12	ng	92
60) 4-Chlorophenyl-phenylether	15.38	204	542986	31.62	ng	# 94
66) 4-Bromophenyl-phenylether	16.27	248	61830	5.04	ng	96
67) Hexachlorobenzene	16.38	284	79939	6.02	ng	93
69) Pentachlorophenol	16.74	266	125962	25.25	ng	95
71) Anthracene	17.21	178	496285	11.99	ng	96
74) Fluoranthene	19.14	202	765172	13.58	ng	95
77) Pyrene	19.50	202	1191755	21.96	ng	96
79) Butylbenzylphthalate	20.41	149	453690	22.80	ng	96
80) Benzo(a)anthracene	21.25	228	1146476	16.88	ng	91
82) Chrysene	21.30	228	841665	12.85	ng	# 88
83) Bis(2-ethylhexyl)phthalate	21.18	149	238233	8.90	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
84) Di-n-octyl phthalate	22.06	149	1258592	30.40	nq	98
85) Indeno(1,2,3-cd)pyrene	25.80	276	1183804	17.68	nq	99
87) Benzo(b)fluoranthene	22.84	252	1715562	27.73	nq	# 98
88) Benzo(k)fluoranthene	22.88	252	795738	13.29	nq	92
89) Benzo(a)pyrene	23.42	252	987643	16.59	nq	97
91) Benzo(a,h,i)perylene	26.51	276	1532320	27.80	nq	# 98

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

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