

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092615\
 Data File : BF081838.D
 Acq On : 26 Sep 2015 20:27
 Operator : UM/IZ
 Sample : G3779-04
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S2-20

Quant Time: Sep 28 07:50:01 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF092415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 25 06:02:49 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.94	152	120657	20.00	ng	-0.01
21) Naphthalene-d8	8.23	136	424029	20.00	ng	-0.01
38) Acenaphthene-d10	9.99	164	208782	20.00	ng	-0.01
63) Phenanthrene-d10	11.49	188	257279	20.00	ng	0.01
75) Chrysene-d12	14.15	240	182005	20.00	ng	0.03
86) Perylene-d12	15.62	264	275229	20.00	ng	0.05

System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	888173	124.53	ng	0.00
7) Phenol-d6	6.56	99	1144354	126.85	ng	-0.01
23) Nitrobenzene-d5	7.51	82	712832	110.92	ng	-0.01
41) 2,4,6-Tribromophenol	10.78	330	217249	119.85	ng	-0.01
44) 2-Fluorobiphenyl	9.30	172	968376	76.14	ng	-0.01
78) Terphenyl-d14	13.06	244	673573	84.54	ng	0.00

Target Compounds				Qvalue		
20) 3+4-Methylphenols	7.34	107	46525	5.58	ng	91
27) 2,4-Dimethylphenol	7.87	122	16632	2.58	ng	# 85
31) Naphthalene	8.26	128	4470995	228.56	ng	# 91
37) 2-Methylnaphthalene	8.95	142	1212256	86.96	ng	# 89
45) 1,1'-Biphenyl	9.41	154	504687	30.92	ng	95
48) Acenaphthylene	9.84	152	357081	16.07	ng	# 94
49) Dimethylphthalate	9.70	163	299875	18.88	ng	# 97
51) Acenaphthene	10.02	154	1252911	97.83	ng	# 88
54) Dibenzofuran	10.19	168	1841506	98.02	ng	# 83
57) Fluorene	10.54	166	1352151	93.79	ng	91
59) Diethylphthalate	10.40	149	74726	4.86	ng	99
70) Phenanthrene	11.54	178	8436437m	573.55	ng	
71) Anthracene	11.58	178	2578781m	181.94	ng	
72) Carbazole	11.73	167	2847665	216.42	ng	# 90
74) Fluoranthene	12.73	202	9407962	634.74	ng	85
77) Pyrene	12.97	202	7701907	561.29	ng	# 85
80) Benzo(a)anthracene	14.14	228	5651419	512.08	ng	# 87
82) Chrysene	14.18	228	2899682	273.96	ng	# 83
85) Indeno(1,2,3-cd)pyrene	17.14	276	2813082	290.31	ng	# 83
87) Benzo(b)fluoranthene	15.21	252	6656300m	407.83	ng	
88) Benzo(k)fluoranthene	15.24	252	1054179m	67.05	ng	
89) Benzo(a)pyrene	15.59	252	3537584	240.06	ng	# 82
90) Dibenzo(a,h)anthracene	17.14	278	755728	49.21	ng	# 77
91) Benzo(g,h,i)perylene	17.59	276	2483669	156.87	ng	# 75

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

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