

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092615\
 Data File : BF081837.D
 Acq On : 26 Sep 2015 19:57
 Operator : UM/IZ
 Sample : G3779-03
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S2-10

Quant Time: Sep 28 08:03:04 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	108039	20.00	ng	-0.01
21) Naphthalene-d8	8.23	136	410794	20.00	ng	-0.01
38) Acenaphthene-d10	9.98	164	185198	20.00	ng	-0.02
63) Phenanthrene-d10	11.47	188	326517	20.00	ng	0.00
75) Chrysene-d12	14.12	240	231627	20.00	ng	0.00
86) Perylene-d12	15.59	264	264571	20.00	ng	0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	795075	124.50	ng	0.00
7) Phenol-d6	6.56	99	994928	123.17	ng	-0.01
23) Nitrobenzene-d5	7.51	82	747899	120.12	ng	-0.01
41) 2,4,6-Tribromophenol	10.78	330	226650	140.96	ng	-0.01
44) 2-Fluorobiphenyl	9.30	172	1014815	89.95	ng	-0.01
78) Terphenyl-d14	13.05	244	715639	70.58	ng	-0.01

Target Compounds				Qvalue		
31) Naphthalene	8.25	128	482466	25.46	ng	99
37) 2-Methylnaphthalene	8.94	142	96646	7.16	ng	# 91
45) 1,1'-Biphenyl	9.41	154	39754	2.75	ng	96
48) Acenaphthylene	9.84	152	63195	3.21	ng	98
49) Dimethylphthalate	9.69	163	302091	21.44	ng	# 97
51) Acenaphthene	10.01	154	203898	17.95	ng	90
54) Dibenzofuran	10.18	168	267226	16.04	ng	# 88
57) Fluorene	10.53	166	226790	17.73	ng	92
59) Diethylphthalate	10.40	149	90663	6.64	ng	97
70) Phenanthrene	11.50	178	2271350m	121.67	ng	
71) Anthracene	11.54	178	659576m	36.67	ng	
72) Carbazole	11.70	167	455776	27.29	ng	96
74) Fluoranthene	12.70	202	2336761	124.23	ng	87
77) Pyrene	12.93	202	1924714	110.22	ng	# 94
80) Benzo(a)anthracene	14.10	228	1097257	78.12	ng	93
82) Chrysene	14.15	228	761542	56.54	ng	91
85) Indeno(1,2,3-cd)pyrene	17.05	276	519814	42.15	ng	# 81
87) Benzo(b)fluoranthene	15.17	252	1218909m	77.69	ng	
88) Benzo(k)fluoranthene	15.19	252	404918m	26.79	ng	
89) Benzo(a)pyrene	15.53	252	785482	55.45	ng	# 84
90) Dibenzo(a,h)anthracene	17.06	278	113719	7.70	ng	# 62
91) Benzo(g,h,i)perylene	17.51	276	462225	30.37	ng	# 74

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

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