

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091415\
 Data File : BF081613.D
 Acq On : 14 Sep 2015 18:14
 Operator : UM/IZ
 Sample : G3597-07 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GP-7(0-5)

Quant Time: Sep 15 03:11:20 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 14 13:51:56 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.23	152	56744	20.00	ng	0.00
21) Naphthalene-d8	11.13	136	222256	20.00	ng	0.00
38) Acenaphthene-d10	15.27	164	100577	20.00	ng	0.00
63) Phenanthrene-d10	18.78	188	174535	20.00	ng	0.01
75) Chrysene-d12	23.40	240	103606	20.00	ng	0.00
86) Perylene-d12	24.84	264	82748	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.38	112	211792	57.91	ng	0.01
7) Phenol-d6	7.65	99	304036	62.80	ng	-0.01
23) Nitrobenzene-d5	9.52	82	194892	42.31	ng	-0.02
41) 2,4,6-Tribromophenol	17.17	330	49622	55.41	ng	-0.01
44) 2-Fluorobiphenyl	13.77	172	322560	41.10	ng	-0.01
78) Terphenyl-d14	22.12	244	186274	41.85	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	14.82	163	64920	8.31	ng	# 97
70) Phenanthrene	18.82	178	36990	3.81	ng	99
74) Fluoranthene	21.44	202	86211	8.21	ng	88
77) Pyrene	21.78	202	107102	13.96	ng	97
80) Benzo(a)anthracene	23.38	228	59945	9.09	ng	95
82) Chrysene	23.42	228	44716	7.56	ng	92
85) Indeno(1,2,3-cd)pyrene	25.89	276	16254	3.75	ng	# 72
87) Benzo(b)fluoranthene	24.49	252	56268m	9.52	ng	
88) Benzo(k)fluoranthene	24.52	252	16934m	3.45	ng	
89) Benzo(a)pyrene	24.78	252	39739	7.94	ng	# 87
91) Benzo(g,h,i)perylene	26.20	276	16962	4.59	ng	# 70

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

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