

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051415\
 Data File : BF079250.D
 Acq On : 15 May 2015 00:17
 Operator : TP/IZ
 Sample : G2177-02
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PS-19(0-3)

Quant Time: May 15 04:03:24 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 15 02:17:41 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.22	152	45903	20.00	ng	0.00
21) Naphthalene-d8	8.80	136	194697	20.00	ng	0.00
38) Acenaphthene-d10	10.97	164	95123	20.00	ng	0.00
63) Phenanthrene-d10	12.81	188	187884	20.00	ng	0.00
75) Chrysene-d12	16.10	240	201709	20.00	ng	-0.14
86) Perylene-d12	17.87	264	217713	20.00	ng	-0.42
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	239710m	89.89	ng	0.00
7) Phenol-d6	6.79	99	339897	96.43	ng	0.00
23) Nitrobenzene-d5	7.91	82	185814	54.85	ng	0.00
41) 2,4,6-Tribromophenol	11.95	330	102117	99.23	ng	0.00
44) 2-Fluorobiphenyl	10.14	172	398844	58.42	ng	-0.01
78) Terphenyl-d14	14.81	244	554381m	65.80	ng	-0.02
Target Compounds						Qvalue
49) Dimethylphthalate	10.65	163	14330	2.05	ng	98
70) Phenanthrene	12.85	178	41866	3.98	ng	98
74) Fluoranthene	14.32	202	140562	12.83	ng	94
77) Pyrene	14.59	202	160354m	13.80	ng	
80) Benzo(a)anthracene	16.09	228	85276	7.72	ng	96
82) Chrysene	16.14	228	63125	5.94	ng	96
85) Indeno(1,2,3-cd)pyrene	19.30	276	54197m	4.00	ng	
87) Benzo(b)fluoranthene	17.42	252	119829m	10.06	ng	
88) Benzo(k)fluoranthene	17.45	252	26008m	2.25	ng	
89) Benzo(a)pyrene	17.81	252	85216	7.77	ng	# 96
91) Benzo(a,h,i)perylene	19.73	276	58689	5.02	ng	# 95

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

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