

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011316\
 Data File : BF084478.D
 Acq On : 14 Jan 2016 9:12
 Operator : SJ/UM
 Sample : H1077-01 5X
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 POST-EX-1-20160108

Quant Time: Jan 14 17:01:38 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF123115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 14 14:15:52 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	320179	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	1320283	20.00	ng	0.00
38) Acenaphthene-d10	9.87	164	541472	20.00	ng	0.00
63) Phenanthrene-d10	11.36	188	903400	20.00	ng	0.00
75) Chrysene-d12	13.99	240	596096	20.00	ng	-0.01
86) Perylene-d12	15.41	264	602819	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	504626	25.49	ng	-0.01
7) Phenol-d6	6.46	99	619998	24.94	ng	-0.01
23) Nitrobenzene-d5	7.39	82	358427	15.11	ng	-0.01
41) 2,4,6-Tribromophenol	10.66	330	119969	21.95	ng	-0.01
44) 2-Fluorobiphenyl	9.20	172	653307	16.10	ng	0.00
78) Terphenyl-d14	12.93	244	372420	13.95	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.60	163	152773	3.95	ng	Qvalue # 92
70) Phenanthrene	11.38	178	224898	4.76	ng	99
74) Fluoranthene	12.57	202	305995	6.20	ng	94
77) Pyrene	12.80	202	279015	5.96	ng	98
80) Benzo(a)anthracene	13.99	228	151254	4.32	ng	95
82) Chrysene	14.02	228	121226m	3.60	ng	
85) Indeno(1,2,3-cd)pyrene	16.79	276	75612	2.84	ng	98
87) Benzo(b)fluoranthene	15.00	252	200492m	5.08	ng	
89) Benzo(a)pyrene	15.35	252	127529	3.81	ng	# 77
91) Benzo(g,h,i)perylene	17.21	276	74899	2.51	ng	95

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

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