

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011316\
 Data File : BF084479.D
 Acq On : 14 Jan 2016 9:40
 Operator : SJ/UM
 Sample : H1077-02 5X
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 POST-EX-2-20160108

Quant Time: Jan 14 17:04:09 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	339350	20.00	ng	0.00
21) Naphthalene-d8	8.12	136	1398863	20.00	ng	0.00
38) Acenaphthene-d10	9.87	164	572856	20.00	ng	0.00
63) Phenanthrene-d10	11.36	188	917714	20.00	ng	0.00
75) Chrysene-d12	13.98	240	638233	20.00	ng	-0.01
86) Perylene-d12	15.41	264	627975	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	332664	15.86	ng	-0.01
7) Phenol-d6	6.46	99	592134	22.47	ng	-0.01
23) Nitrobenzene-d5	7.39	82	359807	14.32	ng	-0.01
41) 2,4,6-Tribromophenol	10.66	330	45189	7.81	ng	-0.01
44) 2-Fluorobiphenyl	9.20	172	658806	15.02	ng	0.00
78) Terphenyl-d14	12.93	244	378454	13.24	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.60	163	135036	3.30	ng	Qvalue # 90
70) Phenanthrene	11.38	178	227296	4.73	ng	98
74) Fluoranthene	12.57	202	455848	9.09	ng	94
77) Pyrene	12.80	202	457930	9.14	ng	98
80) Benzo(a)anthracene	13.98	228	252830	6.75	ng	96
82) Chrysene	14.02	228	198800m	5.52	ng	
85) Indeno(1,2,3-cd)pyrene	16.79	276	116189	4.07	ng	94
87) Benzo(b)fluoranthene	15.00	252	323191m	7.87	ng	
89) Benzo(a)pyrene	15.35	252	213156	6.12	ng	# 88
91) Benzo(g,h,i)perylene	17.21	276	106518	3.43	ng	97

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

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