

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092718\
 Data File : BF109400.D
 Acq On : 27 Sep 2018 21:02
 Operator : JU/SJ
 Sample : J5077-05
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PARCEL-193-092418-1

Manual Integrations
 APPROVED

Sohil
 9/28/2018 2:40:05 PM

Quant Time: Sep 28 05:41:44 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 22 13:32:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	112244	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	429488	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	190310	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	305953	20.00	ng	0.00
75) Chrysene-d12	13.84	240	262713	20.00	ng	0.00
86) Perylene-d12	15.24	264	202174	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	487854	71.59	ng	0.00
7) Phenol-d6	6.35	99	623225	71.67	ng	-0.01
23) Nitrobenzene-d5	7.26	82	400685	55.07	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	134143	71.50	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	734536	58.21	ng	-0.01
78) Terphenyl-d14	12.79	244	555102	51.86	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.46	163	282743	20.427	ng	99
70) Phenanthrene	11.23	178	33797	2.212	ng	98
74) Fluoranthene	12.42	202	63381	3.882	ng	98
77) Pyrene	12.64	202	66302	3.521	ng	97
80) Benzo(a)anthracene	13.83	228	32575	2.059	ng	93
82) Chrysene	13.86	228	34862	2.223	ng	98
87) Benzo(b)fluoranthene	14.84	252	40782m	3.671	ng	
89) Benzo(a)pyrene	15.18	252	22421	2.240	ng	# 90

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

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