

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092718\
 Data File : BF109397.D
 Acq On : 27 Sep 2018 19:43
 Operator : JU/SJ
 Sample : J5077-02
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PARCEL-186-092418-2

Manual Integrations
 APPROVED

Sohil
 9/28/2018 2:40:02 PM

Quant Time: Sep 28 05:32:07 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	114672	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	443327	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	205600	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	345966	20.00	ng	0.00
75) Chrysene-d12	13.84	240	276025	20.00	ng	0.00
86) Perylene-d12	15.24	264	261117	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	503310	72.29	ng	0.00
7) Phenol-d6	6.35	99	642689	72.34	ng	-0.01
23) Nitrobenzene-d5	7.26	82	420659	56.01	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	139990	69.07	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	639953	46.94	ng	-0.01
78) Terphenyl-d14	12.79	244	490559	43.62	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.46	163	297857	19.918	ng	99
74) Fluoranthene	12.42	202	59637	3.231	ng	99
77) Pyrene	12.65	202	52469	2.652	ng	99
87) Benzo(b)fluoranthene	14.84	252	46385m	3.233	ng	
89) Benzo(a)pyrene	15.18	252	28777	2.226	ng	96

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

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