

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092218\
 Data File : BF109237.D
 Acq On : 22 Sep 2018 22:22
 Operator : JU/SJ
 Sample : J4777-19 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-03-092018-A

Manual Integrations
 APPROVED

Sohil
 9/24/2018 11:39:06 AM

Quant Time: Sep 24 07:57:28 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	117926	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	442888	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	187935	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	286964	20.00	ng	0.00
75) Chrysene-d12	13.83	240	235785	20.00	ng	-0.01
86) Perylene-d12	15.23	264	198605	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	397708	55.55	ng	0.00
7) Phenol-d6	6.34	99	505233	55.30	ng	-0.02
23) Nitrobenzene-d5	7.26	82	305815	40.76	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	88994	48.04	ng	-0.02
44) 2-Fluorobiphenyl	9.06	172	557154	44.71	ng	-0.01
78) Terphenyl-d14	12.79	244	367903	38.29	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.46	163	156597	11.456	ng	99
74) Fluoranthene	12.42	202	54548	3.562	ng	96
77) Pyrene	12.65	202	49505	2.930	ng	98
82) Chrysene	13.86	228	29403	2.089	ng	99
87) Benzo(b)fluoranthene	14.84	252	34573m	3.168	ng	
89) Benzo(a)pyrene	15.17	252	21785	2.215	ng	# 92

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

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