

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091018\
 Data File : BF108891.D
 Acq On : 10 Sep 2018 14:29
 Operator : JU/SJ
 Sample : J4776-01 2X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-02-090718-A

Manual Integrations
 APPROVED

Sohil
 9/11/2018 12:01:30 PM

Quant Time: Sep 11 09:01:14 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF090518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 05 18:03:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.74	152	205991	20.00	ng	0.00
21) Naphthalene-d8	8.02	136	749562	20.00	ng	0.00
38) Acenaphthene-d10	9.77	164	315276	20.00	ng	0.00
63) Phenanthrene-d10	11.24	188	530023	20.00	ng	0.00
75) Chrysene-d12	13.87	240	516743	20.00	ng	0.00
86) Perylene-d12	15.28	264	442560	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.34	112	582415	46.84	ng	0.00
7) Phenol-d6	6.37	99	722877	45.18	ng	0.00
23) Nitrobenzene-d5	7.29	82	435794	36.38	ng	-0.01
41) 2,4,6-Tribromophenol	10.55	330	133501	44.36	ng	0.00
44) 2-Fluorobiphenyl	9.10	172	776812	42.03	ng	0.00
78) Terphenyl-d14	12.83	244	642753	29.01	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.48	163	197709	8.966	ng	Qvalue 100
70) Phenanthrene	11.27	178	80064	3.163	ng	99
74) Fluoranthene	12.45	202	154433	6.162	ng	96
77) Pyrene	12.68	202	165824	4.170	ng	98
80) Benzo(a)anthracene	13.86	228	100040	3.020	ng	94
82) Chrysene	13.89	228	92605m	2.888	ng	
87) Benzo(b)fluoranthene	14.88	252	114448m	4.756	ng	
89) Benzo(a)pyrene	15.22	252	67537	3.110	ng	# 83

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

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