

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071218\
 Data File : BF107347.D
 Acq On : 12 Jul 2018 18:54
 Operator : JU/SJ
 Sample : J3832-01 2X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-05-071018

Manual Integrations
 APPROVED

Sohil
 7/13/2018 11:17:14 AM

Quant Time: Jul 13 03:29:20 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 11 16:39:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	60912	20.00	ng	0.00
21) Naphthalene-d8	8.21	136	218021	20.00	ng	0.00
38) Acenaphthene-d10	9.98	164	94126	20.00	ng	0.00
63) Phenanthrene-d10	11.48	188	181275	20.00	ng	0.00
75) Chrysene-d12	14.14	240	181892	20.00	ng	0.00
86) Perylene-d12	15.68	264	124494	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	227079	63.13	ng	0.00
7) Phenol-d6	6.58	99	274140	61.03	ng	0.00
23) Nitrobenzene-d5	7.50	82	164951	42.05	ng	0.00
41) 2,4,6-Tribromophenol	10.78	330	62861	57.62	ng	0.00
44) 2-Fluorobiphenyl	9.29	172	281225	43.34	ng	0.00
78) Terphenyl-d14	13.07	244	328914	43.80	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.68	163	46628	6.605	ng	98
70) Phenanthrene	11.50	178	18377	2.102	ng	98
74) Fluoranthene	12.70	202	33194	3.444	ng	97
77) Pyrene	12.94	202	31520	2.628	ng	98
87) Benzo(b)fluoranthene	15.22	252	17303m	2.237	ng	

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

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