

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062918\
 Data File : BF106970.D
 Acq On : 29 Jun 2018 13:59
 Operator : JU/SJ
 Sample : J3709-05
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EME-FW-TS005-01

Manual Integrations
 APPROVED

Sohil
 6/30/2018 10:57:57 AM

Quant Time: Jun 29 19:00:00 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 29 11:40:02 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	71105	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	269343	20.00	ng	-0.01
38) Acenaphthene-d10	10.00	164	130846	20.00	ng	-0.01
63) Phenanthrene-d10	11.49	188	225912	20.00	ng	-0.01
75) Chrysene-d12	14.14	240	205222	20.00	ng	-0.02
86) Perylene-d12	15.69	264	168783	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	409663	97.56	ng	0.00
7) Phenol-d6	6.60	99	522438	99.63	ng	-0.02
23) Nitrobenzene-d5	7.52	82	324789	67.01	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	153766	101.39	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	559717	69.76	ng	-0.02
78) Terphenyl-d14	13.07	244	511544	60.38	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.70	163	73711	7.511	ng	100
74) Fluoranthene	12.71	202	48888	4.071	ng	93
77) Pyrene	12.94	202	43078	3.183	ng	98
80) Benzo(a)anthracene	14.13	228	28516	2.280	ng	96
82) Chrysene	14.17	228	28078m	2.440	ng	
87) Benzo(b)fluoranthene	15.23	252	39256m	3.744	ng	
89) Benzo(a)pyrene	15.62	252	22389	2.402	ng	# 93

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

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