

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062718\
 Data File : BF106885.D
 Acq On : 27 Jun 2018 21:06
 Operator : JU/SJ
 Sample : J3242-11 2X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-062618-F

Manual Integrations
 APPROVED

Sohil
 6/28/2018 2:21:37 PM

Quant Time: Jun 28 04:21:17 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 25 18:50:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	21373	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	81379	20.00	ng	-0.01
38) Acenaphthene-d10	10.00	164	36737	20.00	ng	-0.01
63) Phenanthrene-d10	11.49	188	63049	20.00	ng	-0.01
75) Chrysene-d12	14.14	240	66619	20.00	ng	-0.02
86) Perylene-d12	15.68	264	48511	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	121868	96.55	ng	-0.01
7) Phenol-d6	6.59	99	155929	98.92	ng	-0.02
23) Nitrobenzene-d5	7.51	82	91356	62.39	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	40771	95.75	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	186879	86.33	ng	-0.02
78) Terphenyl-d14	13.07	244	194327	70.66	ng	-0.01
Target Compounds						
10) Phenol	6.60	94	4104	2.281	ng	# 63
49) Dimethylphthalate	9.70	163	37208	13.504	ng	# 98
74) Fluoranthene	12.71	202	7242	2.161	ng	94
77) Pyrene	12.94	202	11123	2.532	ng	97
87) Benzo(b)fluoranthene	15.22	252	6778m	2.249	ng	

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

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