

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF063018\
 Data File : BF107026.D
 Acq On : 30 Jun 2018 20:24
 Operator : JU/SJ
 Sample : J3638-02 2X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P-2

Manual Integrations
 APPROVED

Sohil
 7/5/2018 4:32:19 PM

Quant Time: Jul 02 07:16:18 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 14:38:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	68734	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	241619	20.00	ng	-0.01
38) Acenaphthene-d10	10.00	164	105477	20.00	ng	-0.01
63) Phenanthrene-d10	11.49	188	213039	20.00	ng	-0.01
75) Chrysene-d12	14.14	240	186502	20.00	ng	-0.02
86) Perylene-d12	15.68	264	133715	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	226812	55.88	ng	0.00
7) Phenol-d6	6.60	99	272470	53.75	ng	-0.02
23) Nitrobenzene-d5	7.51	82	150382	34.59	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	71834	58.76	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	310574	42.46	ng	-0.02
78) Terphenyl-d14	13.07	244	393660	51.13	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.70	163	36839	4.657	ng	97
70) Phenanthrene	11.51	178	69042	6.719	ng	97
74) Fluoranthene	12.71	202	182022	16.072	ng	94
77) Pyrene	12.94	202	168774	13.723	ng	99
80) Benzo(a)anthracene	14.13	228	104321	9.178	ng	98
82) Chrysene	14.17	228	99670	9.531	ng	98
85) Indeno(1,2,3-cd)pyrene	17.25	276	69025	7.778	ng	# 87
87) Benzo(b)fluoranthene	15.22	252	132999m	16.012	ng	
88) Benzo(k)fluoranthene	15.25	252	45837m	5.795	ng	
89) Benzo(a)pyrene	15.61	252	86795	11.754	ng	96
91) Benzo(g,h,i)perylene	17.74	276	70908	11.593	ng	# 90

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

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