

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF063018\
 Data File : BF107010.D
 Acq On : 30 Jun 2018 13:16
 Operator : JU/SJ
 Sample : J3248-09 5X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-03-062818-B

Manual Integrations
 APPROVED

Sohil
 7/2/2018 12:29:56 PM

Quant Time: Jul 02 05:05:34 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	63823	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	228879	20.00	ng	-0.01
38) Acenaphthene-d10	10.00	164	95709	20.00	ng	-0.01
63) Phenanthrene-d10	11.49	188	176571	20.00	ng	-0.01
75) Chrysene-d12	14.14	240	216174	20.00	ng	-0.02
86) Perylene-d12	15.69	264	173555	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	96902	25.71	ng	-0.01
7) Phenol-d6	6.60	99	113481	24.11	ng	-0.02
23) Nitrobenzene-d5	7.51	82	63886	15.51	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	24887	22.44	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	113333	6.40	ng	-0.02
78) Terphenyl-d14	13.07	244	134338	15.05	ng	-0.01
Target Compounds						
70) Phenanthrene	11.51	178	65604	7.703	ng	99
71) Anthracene	11.57	178	27490	3.162	ng	99
74) Fluoranthene	12.71	202	104025	11.082	ng	95
77) Pyrene	12.94	202	141597	9.933	ng	99
80) Benzo(a)anthracene	14.14	228	73034	5.543	ng	92
82) Chrysene	14.17	228	72317m	5.966	ng	
85) Indeno(1,2,3-cd)pyrene	17.27	276	22012	2.140	ng	# 87
87) Benzo(b)fluoranthene	15.23	252	64351m	5.969	ng	
89) Benzo(a)pyrene	15.62	252	52105	5.437	ng	94
91) Benzo(g,h,i)perylene	17.75	276	23032m	2.901	ng	

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

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