

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF063018\
 Data File : BF107011.D
 Acq On : 30 Jun 2018 13:42
 Operator : JU/SJ
 Sample : J3248-11 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 7/2/2018 12:29:59 PM

Quant Time: Jul 02 05:15:02 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	57531	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	202865	20.00	ng	-0.01
38) Acenaphthene-d10	10.00	164	85209	20.00	ng	-0.01
63) Phenanthrene-d10	11.49	188	171492	20.00	ng	-0.01
75) Chrysene-d12	14.14	240	212783	20.00	ng	-0.02
86) Perylene-d12	15.69	264	157533	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	77526	22.82	ng	0.00
7) Phenol-d6	6.60	99	90807	21.40	ng	-0.02
23) Nitrobenzene-d5	7.51	82	50211	13.75	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	21179	21.45	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	90760	3.97	ng	-0.02
78) Terphenyl-d14	13.07	244	121059	13.78	ng	-0.02
Target Compounds						
70) Phenanthrene	11.51	178	144055	17.416	ng	98
71) Anthracene	11.57	178	33414	3.958	ng	100
74) Fluoranthene	12.71	202	199926	21.930	ng	95
77) Pyrene	12.94	202	192379	13.710	ng	98
80) Benzo(a)anthracene	14.13	228	100526	7.752	ng	98
82) Chrysene	14.17	228	92605m	7.762	ng	
85) Indeno(1,2,3-cd)pyrene	17.27	276	25587	2.527	ng	# 92
87) Benzo(b)fluoranthene	15.23	252	86634m	8.853	ng	
88) Benzo(k)fluoranthene	15.26	252	22289m	2.392	ng	
89) Benzo(a)pyrene	15.62	252	56784	6.527	ng	98
91) Benzo(g,h,i)perylene	17.75	276	26454	3.671	ng	# 91

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

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