

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070618\
 Data File : BF107185.D
 Acq On : 6 Jul 2018 20:46
 Operator : JU/SJ
 Sample : J3868-07
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-7-D

Manual Integrations
 APPROVED

Sohil
 7/9/2018 1:58:41 PM

Quant Time: Jul 07 00:42:37 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	6.94	152	61978	20.00	ng	-0.02
21) Naphthalene-d8	8.23	136	231831	20.00	ng	-0.02
38) Acenaphthene-d10	9.98	164	101005	20.00	ng	-0.02
63) Phenanthrene-d10	11.48	188	164252	20.00	ng	-0.02
75) Chrysene-d12	14.14	240	171798	20.00	ng	-0.03
86) Perylene-d12	15.67	264	131837	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	335943	91.78	ng	-0.01
7) Phenol-d6	6.59	99	423358	92.62	ng	-0.02
23) Nitrobenzene-d5	7.51	82	261502	62.69	ng	-0.02
41) 2,4,6-Tribromophenol	10.78	330	95835	81.86	ng	-0.02
44) 2-Fluorobiphenyl	9.30	172	454694	74.35	ng	-0.02
78) Terphenyl-d14	13.06	244	428728	60.45	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.70	163	131286	17.330	ng	Qvalue # 99
70) Phenanthrene	11.51	178	119867	15.131	ng	97
71) Anthracene	11.56	178	44978	5.562	ng	96
74) Fluoranthene	12.70	202	284142	32.541	ng	95
77) Pyrene	12.93	202	276378	24.396	ng	99
80) Benzo(a)anthracene	14.12	228	154730	14.778	ng	99
82) Chrysene	14.16	228	135028m	14.017	ng	
85) Indeno(1,2,3-cd)pyrene	17.24	276	72237	8.837	ng	# 88
87) Benzo(b)fluoranthene	15.21	252	184290m	22.503	ng	
88) Benzo(k)fluoranthene	15.24	252	43568m	5.586	ng	
89) Benzo(a)pyrene	15.60	252	126420m	17.364	ng	
90) Dibenzo(a,h)anthracene	17.24	278	17177m	2.784	ng	
91) Benzo(a,h,i)perylene	17.72	276	76822	12.738	ng	# 90

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

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