

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF041919\
 Data File : BF113867.D
 Acq On : 19 Apr 2019 19:05
 Operator : JU/SJ
 Sample : K2199-15 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-03-041819-B

Manual Integrations
 APPROVED

Sohil
 4/22/2019 7:27:04 PM

Quant Time: Apr 20 05:00:01 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF041819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 18 13:55:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	163674	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	558462	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	253416	20.00	ng	0.00
64) Phenanthrene-d10	11.43	188	512830	20.00	ng	0.02
76) Chrysene-d12	14.35	240	447389m	20.00	ng	0.31
87) Perylene-d12	16.17	264	351064m	20.00	ng	0.66
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	211914	21.53	ng	0.00
7) Phenol-d6	6.52	99	269328	21.94	ng	0.00
23) Nitrobenzene-d5	7.45	82	130237	14.20	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	63330	22.62	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	258897	15.58	ng	0.00
79) Terphenyl-d14	13.10	244	332242	15.57	ng	0.10
Target Compounds						
71) Phenanthrene	11.45	178	165543	6.338	ng	97
75) Fluoranthene	12.69	202	279891	10.111	ng	97
78) Pyrene	12.94	202	258031	8.184	ng	97
83) Chrysene	14.34	228	123927m	4.774	ng	
86) Indeno(1,2,3-cd)pyrene	17.72	276	42687m	2.014	ng	
88) Benzo(b)fluoranthene	15.69	252	103155m	4.713	ng	
90) Benzo(a)pyrene	16.10	252	70588m	3.648	ng	
92) Benzo(g,h,i)perylene	18.18	276	39197m	2.240	ng	

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

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