

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121018\
 Data File : BF111262.D
 Acq On : 11 Dec 2018 00:26
 Operator : JU/SJ
 Sample : J6234-01 5X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 VNJ-218

Manual Integrations
 APPROVED

Sohil
 12/11/2018 2:23:45 PM

Quant Time: Dec 11 06:19:05 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 10 01:15:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	517515	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	1892599	20.00	ng	0.00
39) Acenaphthene-d10	9.84	164	542578	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	930316	20.00	ng	0.00
76) Chrysene-d12	13.96	240	638167	20.00	ng	0.00
87) Perylene-d12	15.40	264	463394	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	701593	21.41	ng	0.00
7) Phenol-d6	6.42	99	886318	21.76	ng	-0.01
23) Nitrobenzene-d5	7.36	82	492985	14.61	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	84003	17.48	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	747496	16.77	ng	0.00
79) Terphenyl-d14	12.90	244	470619	17.69	ng	0.00
Target Compounds						
71) Phenanthrene	11.34	178	133806	2.852	ng	93
75) Fluoranthene	12.53	202	191057	4.457	ng	98
78) Pyrene	12.76	202	180661	3.962	ng	96
81) Benzo(a)anthracene	13.94	228	80752	2.268	ng	94
83) Chrysene	13.98	228	76753	2.091	ng	97
88) Benzo(b)fluoranthene	14.98	252	94037m	3.669	ng	
90) Benzo(a)pyrene	15.33	252	58148	2.464	ng	# 80
92) Benzo(g,h,i)perylene	17.23	276	40323	2.173	ng	# 90

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

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