

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF041919\
 Data File : BF113871.D
 Acq On : 19 Apr 2019 20:48
 Operator : JU/SJ
 Sample : K2201-09 5X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-01-041819-E

Manual Integrations
 APPROVED

Sohil
 4/22/2019 7:27:15 PM

Quant Time: Apr 20 05:12:07 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	163957	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	563868	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	258613	20.00	ng	0.00
64) Phenanthrene-d10	11.43	188	511615	20.00	ng	0.02
76) Chrysene-d12	14.41	240	422573m	20.00	ng	0.36
87) Perylene-d12	16.26	264	336487m	20.00	ng	0.75
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	207867	21.08	ng	0.00
7) Phenol-d6	6.52	99	264203	21.49	ng	0.00
23) Nitrobenzene-d5	7.45	82	123542	13.34	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	62475	21.87	ng	0.01
45) 2-Fluorobiphenyl	9.26	172	271790	16.03	ng	0.00
79) Terphenyl-d14	13.13	244	340872	16.92	ng	0.13
Target Compounds						
71) Phenanthrene	11.46	178	59852	2.297	ng	98
75) Fluoranthene	12.71	202	129997	4.707	ng	97
78) Pyrene	12.96	202	128618	4.319	ng	95
81) Benzo(a)anthracene	14.39	228	59056m	2.391	ng	
83) Chrysene	14.44	228	58013m	2.366	ng	

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

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