

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF010819\
 Data File : BF111907.D
 Acq On : 8 Jan 2019 15:20
 Operator : JU/SJ
 Sample : K1012-03 2X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-02-010419-B

Manual Integrations
 APPROVED

Sohil
 1/9/2019 10:36:42 AM

Quant Time: Jan 08 18:09:00 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF010719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 07 14:33:51 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.88	152	148544	20.00	ng	-0.01
21) Naphthalene-d8	8.16	136	517444	20.00	ng	-0.01
39) Acenaphthene-d10	9.92	164	197220	20.00	ng	-0.02
64) Phenanthrene-d10	11.40	188	374181	20.00	ng	-0.01
76) Chrysene-d12	14.04	240	430124	20.00	ng	-0.01
87) Perylene-d12	15.53	264	374265	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	479238	55.19	ng	0.00
7) Phenol-d6	6.52	99	570464	50.98	ng	-0.01
23) Nitrobenzene-d5	7.44	82	379316	39.18	ng	-0.02
42) 2,4,6-Tribromophenol	10.70	330	122825	47.88	ng	-0.02
45) 2-Fluorobiphenyl	9.23	172	491741	36.36	ng	-0.02
79) Terphenyl-d14	12.99	244	534381	23.58	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.63	163	35970	2.428	ng	Qvalue # 97
71) Phenanthrene	11.43	178	148845	7.440	ng	99
75) Fluoranthene	12.62	202	268972	13.097	ng	98
78) Pyrene	12.85	202	272877	9.220	ng	98
81) Benzo(a)anthracene	14.03	228	145597	5.882	ng	98
83) Chrysene	14.07	228	128971m	5.464	ng	
86) Indeno(1,2,3-cd)pyrene	17.02	276	39710	2.308	ng	96
88) Benzo(b)fluoranthene	15.09	252	143791m	6.423	ng	
90) Benzo(a)pyrene	15.46	252	92769	4.713	ng	# 90
92) Benzo(g,h,i)perylene	17.46	276	33289	2.150	ng	# 91

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

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