

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF011819\
 Data File : BF112148.D
 Acq On : 19 Jan 2019 2:02
 Operator : JU/SJ
 Sample : K1015-09 5X
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-01-011719-A

Manual Integrations
 APPROVED

Sohil
 1/21/2019 12:01:04 PM

Quant Time: Jan 19 03:58:32 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF011619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 19 01:02:12 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.86	152	246124	20.00	ng	0.00
21) Naphthalene-d8	8.14	136	940736	20.00	ng	0.00
39) Acenaphthene-d10	9.89	164	406251	20.00	ng	0.00
64) Phenanthrene-d10	11.38	188	676515	20.00	ng	0.00
76) Chrysene-d12	14.02	240	568495	20.00	ng	0.00
87) Perylene-d12	15.50	264	442067	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	282738	20.85	ng	-0.01
7) Phenol-d6	6.49	99	363246	21.41	ng	-0.01
23) Nitrobenzene-d5	7.41	82	204081	14.99	ng	-0.02
42) 2,4,6-Tribromophenol	10.68	330	70613	16.11	ng	-0.01
45) 2-Fluorobiphenyl	9.21	172	432963	15.63	ng	0.00
79) Terphenyl-d14	12.96	244	368843	13.80	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.60	163	111628	3.768	ng	98
71) Phenanthrene	11.40	178	423922	12.451	ng	96
72) Anthracene	11.45	178	171066	4.960	ng	97
75) Fluoranthene	12.59	202	751423	20.904	ng	100
78) Pyrene	12.82	202	670250	17.870	ng	99
81) Benzo(a)anthracene	14.01	228	368472	11.304	ng	99
83) Chrysene	14.04	228	309014	10.019	ng	97
86) Indeno(1,2,3-cd)pyrene	16.97	276	121163	4.258	ng	93
88) Benzo(b)fluoranthene	15.06	252	351218m	13.996	ng	
89) Benzo(k)fluoranthene	15.09	252	129428m	5.329	ng	
90) Benzo(a)pyrene	15.43	252	236727	10.287	ng	96
92) Benzo(a,h,i)perylene	17.42	276	111717	5.548	ng	98

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

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