

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF012519\
 Data File : BF112258.D
 Acq On : 26 Jan 2019 3:20
 Operator : JU/SJ
 Sample : K1223-09RE 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TS-URE

Manual Integrations
 APPROVED

mohammad
 1/29/2019 12:32:00 AM

Quant Time: Jan 28 01:03:53 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF012519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 25 16:47:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	278624	20.00	ng	-0.01
21) Naphthalene-d8	8.12	136	1074998	20.00	ng	-0.01
39) Acenaphthene-d10	9.87	164	488473	20.00	ng	-0.01
64) Phenanthrene-d10	11.36	188	784998	20.00	ng	-0.01
76) Chrysene-d12	14.00	240	529905	20.00	ng	-0.01
87) Perylene-d12	15.47	264	520967	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	746794	56.93	ng	0.00
7) Phenol-d6	6.49	99	930162	51.03	ng	-0.02
23) Nitrobenzene-d5	7.40	82	574309	30.78	ng	-0.02
42) 2,4,6-Tribromophenol	10.67	330	225848	39.72	ng	-0.02
45) 2-Fluorobiphenyl	9.20	172	1046976	30.31	ng	-0.01
79) Terphenyl-d14	12.94	244	697648	24.80	ng	-0.02
Target Compounds						
50) Dimethylphthalate	9.59	163	105996	2.793	ng	98
71) Phenanthrene	11.39	178	442693	10.847	ng	98
72) Anthracene	11.44	178	85990	2.115	ng	99
75) Fluoranthene	12.57	202	502749	11.486	ng	99
78) Pyrene	12.80	202	432138	11.779	ng	97
81) Benzo(a)anthracene	13.99	228	207803	6.671	ng	99
83) Chrysene	14.03	228	192152	6.462	ng	96
84) Bis(2-ethylhexyl)phthalate	13.97	149	59568	2.364	ng	100
86) Indeno(1,2,3-cd)pyrene	16.94	276	83960	3.563	ng	95
88) Benzo(b)fluoranthene	15.04	252	241491m	7.751	ng	
89) Benzo(k)fluoranthene	15.07	252	94897m	3.180	ng	
90) Benzo(a)pyrene	15.41	252	159973	5.706	ng	# 94
92) Benzo(a,h,i)perylene	17.39	276	79458	3.728	ng	95

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

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